```
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PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2


* * * * * * * * * *   Welcome to STN International   * * * * * * * * * *

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 NEWS  4  DEC 14   2006 MeSH terms loaded in MEDLINE/LMEDLINE
 NEWS  5  DEC 14   2006 MeSH terms loaded for MEDLINE file segment of TOXCENTER
 NEWS  6  DEC 14   CA/CAplus to be enhanced with updated IPC codes
 NEWS  7  DEC 21   IPC search and display fields enhanced in CA/CAplus with the
                   IPC reform
 NEWS  8  DEC 23   New IPC8 SEARCH, DISPLAY, and SELECT fields in USPATFULL/
                   USPAT2
 NEWS  9  JAN 13   IPC 8 searching in IFIPAT, IFIUDB, and IFICDB
 NEWS 10  JAN 13   New IPC 8 SEARCH, DISPLAY, and SELECT enhancements added to
                   INPADOC
 NEWS 11  JAN 17   Pre-1988 INPI data added to MARPAT
 NEWS 12  JAN 17   IPC 8 in the WPI family of databases including WPIFV
 NEWS 13  JAN 30   Saved answer limit increased
 NEWS 14  JAN 31   Monthly current-awareness alert (SDI) frequency
                   added to TULSA
 NEWS 15  FEB 21   STN AnaVist, Version 1.1, lets you share your STN AnaVist
                   visualization results
 NEWS 16  FEB 22   Status of current WO (PCT) information on STN
 NEWS 17  FEB 22   The IPC thesaurus added to additional patent databases on STN
 NEWS 18  FEB 22   Updates in EPFULL; IPC 8 enhancements added
 NEWS 19  FEB 27   New STN AnaVist pricing effective March 1, 2006
 NEWS 20  FEB 28   MEDLINE/LMEDLINE reload improves functionality
 NEWS 21  FEB 28   TOXCENTER reloaded with enhancements
 NEWS 22  FEB 28   REGISTRY/ZREGISTRY enhanced with more experimental spectral
                   property data
 NEWS 23  MAR 01   INSPEC reloaded and enhanced
 NEWS 24  MAR 03   Updates in PATDPA; addition of IPC 8 data without attributes
 NEWS 25  MAR 08   X.25 communication option no longer available after June 2006

 NEWS EXPRESS  FEBRUARY 15 CURRENT VERSION FOR WINDOWS IS V8.01a,
               CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
               AND CURRENT DISCOVER FILE IS DATED 19 DECEMBER 2005.
               V8.0 AND V8.01 USERS CAN OBTAIN THE UPGRADE TO V8.01a AT
               http://download.cas.org/express/v8.0-Discover/

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* * * * * * * * * * * * * * STN Columbus * * * * * * * * * * * * * * * *

FILE 'HOME' ENTERED AT 08:34:46 ON 09 MAR 2006

=> fil caplus
COST IN U.S. DOLLARS

	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'CAPLUS' ENTERED AT 08:35:04 ON 09 MAR 2006
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FILE COVERS 1907 - 9 Mar 2006 VOL 144 ISS 11
FILE LAST UPDATED: 8 Mar 2006 (20060308/ED)

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They are available for your review at:

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=> s jp200017145/pn
L1 0 JP200017145/PN

=> s jp0017145/pn
L2 0 JP0017145/PN

=> s jp2000-17145/pn
L3 1 JP2000-17145/PN
(JP2000017145/PN)

=> d

L3 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN
AN 2000:37940 CAPLUS
DN 132:79372
TI Thermosetting novolak resin compositions with good curability, and molding
materials therefrom
IN Oka, Wataru; Orihara, Tamotsu
PA Sumitomo Bakelite Co., Ltd., Japan
SO Jpn. Kokai Tokkyo Koho, 6 pp.
CODEN: JKXXAF
DT Patent
LA Japanese
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 2000017145	A2	20000118	JP 1998-184487	19980630
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PRAI	JP 1998-184487		19980630		

OS MARPAT 132:79372

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=> d ibib abs
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L3 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2000:37940 CAPLUS
DOCUMENT NUMBER: 132:79372
TITLE: Thermosetting novolak resin compositions with good curability, and molding materials therefrom
INVENTOR(S): Oka, Wataru; Orihara, Tamotsu
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 6 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2000017145	A2	20000118	JP 1998-184487	19980630

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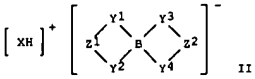
PRIORITY APPLN. INFO.: JP 1998-184487 19980630

OTHER SOURCE(S): MARPAT 132:79372
GI



II

AB The compns. contain novolak resins, hexamethylenetetramine (I), and onium borates II (X = N-containing heterocyclic compound; Z1, Z2 = aromatic or alicyclic group; Y1-Y4 = proton donor group residue). Thus, novolak resin 100, I 16, II (X = DBU, Y1Z1Y2 = Y3Z2Y4 = O-o-C6H4CO2) 8 parts, and fillers are mixed and transfer-molded to give a test piece showing Barcol hardness 68 and bending strength 102 and 63 MPa, at room temperature and 120°, resp.

```
=> s jp11-209583/pn
L4            1 JP11-209583/PN
                (JP11209583/PN)

=> d ibib abs
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L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1999:481505 CAPLUS
DOCUMENT NUMBER: 131:130718
TITLE: Epoxy resin compositions for prepregs and laminated circuit boards
INVENTOR(S): Go, Yoshiyuki; Miyake, Sumiya; Nagata, Hiroshi; Okubo, Akiko; Kobayashi, Minoru
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 8 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 11209583	A2	19990803	JP 1998-11829	19980123

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PRIORITY APPLN. INFO.: JP 1998-11829 19980123

OTHER SOURCE(S): MARPAT 131:130718

AB Title epoxy resin compns., which is stable at room temperature and fat curing upon heating, comprise an epoxy resin, a polyamine curing agent, and a curing accelerator of an onium borate. Thus a bisphenol A-based epoxy resin 100 parts, diaminodiphenylmethane 0.2 parts, dicyandiamide 0.3 parts, and DBU salt of 2-hydroxybenzoic acid boron complex 2.5 parts were mixed to give an epoxy composition of this invention. Two pieces of glass cloths were impregnated with the above composition to two prepregs which were laminated with an printed circuit board by vacuum compression while heating to give a laminated board.

```
=> s jp11-171981/pn
L5            1 JP11-171981/PN
                (JP11171981/PN)

=> d ibib abs
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L5 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1999:407164 CAPLUS
DOCUMENT NUMBER: 131:103041
TITLE: Storage-stable epoxy resin compositions containing ammonium borates as latent crosslinking accelerators
INVENTOR(S): Miyake, Sumiya; Go, Yoshiyuki; Nagata, Hiroshi; Okubo, Akiko; Kobayashi, Minoru
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 8 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 11171981	A2	19990629	JP 1997-340126	19971210
JP 3690710	B2	20050831		
PRIORITY APPLN. INFO.:			JP 1997-340126	19971210

OTHER SOURCE(S): MARPAT 131:103041
GI For diagram(s), see printed CA Issue.
AB Title compns., useful for elec. an electronic devices, etc., contain hardeners and X+ BY1Y2Y3Y4- [X+ = (substituted) ammonium; ≥1 of Y1-Y4 = H+-donating group residue after releasing 1 H+, the rest of Y1-Y4 = aromatic, heterocyclic, or aliphatic group] or I (Y9-Y10 are same as Y1-Y4; Y11-Y12 = H+-donating group residue after release of H+). Thus, o-cresol novolak epoxy resin (EOCN 102065) 67, phenol novolak 33, pulverized fused silica 300, carnauba wax 2, and Ph4N+ (BzO)4B- 3.1 part was mixed and roll-kneaded at 90° for 5 min to give title composition having initial spiral flow 83 cm and 79 cm after 3-day storage at 40°.

```
=> fil reg
COST IN U.S. DOLLARS                                    SINCE FILE      TOTAL
                                                           ENTRY    SESSION
FULL ESTIMATED COST                                        23.25      23.46

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)              SINCE FILE      TOTAL
                                                           ENTRY    SESSION
CA SUBSCRIBER PRICE                                        -2.25      -2.25

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DICTIONARY FILE UPDATES:    8 MAR 2006  HIGHEST RN 876273-86-8

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TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

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**********************************************************************
*                                                                    *
* The CA roles and document type information have been removed from *
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* available and contains the CA role and document type information.  *
*                                                                    *
**********************************************************************

Structure search iteration limits have been increased.  See HELP SLIMITS
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REGISTRY includes numerically searchable data for experimental and
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experimental property data in the original document.  For information
on property searching in REGISTRY, refer to:

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=> fil stnguide
COST IN U.S. DOLLARS                                    SINCE FILE      TOTAL
                                                           ENTRY    SESSION
FULL ESTIMATED COST                                         0.44      23.90

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)              SINCE FILE      TOTAL
                                                           ENTRY    SESSION
CA SUBSCRIBER PRICE                                         0.00      -2.25

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AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE
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=> log y
COST IN U.S. DOLLARS                                    SINCE FILE      TOTAL
                                                          ENTRY    SESSION
FULL ESTIMATED COST                                        15.79     404.57

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)              SINCE FILE      TOTAL
                                                          ENTRY    SESSION
CA SUBSCRIBER PRICE                                        -2.25      -4.50

STN INTERNATIONAL LOGOFF AT 08:49:36 ON 09 MAR 2006


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TERMINAL (ENTER 1, 2, 3, OR ?):2


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 NEWS  5  DEC 14   2006 MeSH terms loaded for MEDLINE file segment of TOXCENTER
 NEWS  6  DEC 14   CA/CAplus to be enhanced with updated IPC codes
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 NEWS  8  DEC 23   New IPC8 SEARCH, DISPLAY, and SELECT fields in USPATFULL/
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 NEWS  9  JAN 13   IPC 8 searching in IFIPAT, IFIUDB, and IFICDB
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                   property data
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               AND CURRENT DISCOVER FILE IS DATED 19 DECEMBER 2005.
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FILE 'HOME' ENTERED AT 08:50:55 ON 09 MAR 2006

=> fil caplus
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                       ENTRY    SESSION
FULL ESTIMATED COST                                     0.21       0.21

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FILE COVERS 1907 - 9 Mar 2006  VOL 144 ISS 11
FILE LAST UPDATED: 8 Mar 2006  (20060308/ED)

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=> s us20020015883
L1            0 US20020015883

=> s us20020015883/pn
L2            1 US20020015883/PN
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ENTER DISPLAY CODE (TI) OR ?:rn
E1 THROUGH E14 ASSIGNED

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FULL ESTIMATED COST                                     4.44       4.65

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DICTIONARY FILE UPDATES:    8 MAR 2006  HIGHEST RN 876273-86-8

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*********************************************************************
*                                                                   *
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added,   *
* effective March 20, 2005.  A new display format, IDERL, is now    *
* available and contains the CA role and document type information. *
*                                                                   *
*********************************************************************

Structure search iteration limits have been increased.  See HELP SLIMITS
for details.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document.  For information
on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

=> s e1-e14
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             1 156762-86-6/BI
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L3 ANSWER 1 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
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ED Entered STN: 19 Dec 2001
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SR CA
LC STN Files: CA, CAPLUS, USPATFULL

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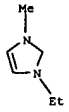

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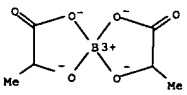

1 REFERENCES IN FILE CA (1907 TO DATE)
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L3 ANSWER 2 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
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ED Entered STN: 19 Dec 2001
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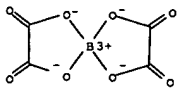

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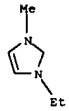

ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE
1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 3 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
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ED Entered STN: 19 Dec 2001
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LC STN Files: CA, CAPLUS, USPATFULL

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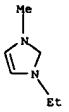

ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

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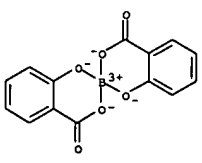
CCI CCS



1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 4 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN **376650-04-3** REGISTRY
ED Entered STN: 19 Dec 2001
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SR CA
LC STN Files: CA, CAPLUS, USPATFULL

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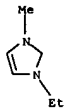
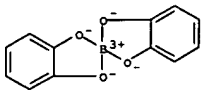
CRN 65039-03-4
CMF C6 H11 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

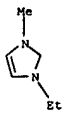
CM 2

CRN 16986-25-7
CMF C12 H8 B O4
CCI CCS



1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

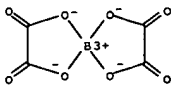
L3 ANSWER 5 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN **250358-46-4** REGISTRY
ED Entered STN: 09 Dec 1999
CN 1H-Imidazolium, 1-ethyl-3-methyl-, hydroxide (9CI) (CA INDEX NAME)
OTHER NAMES:
CN 1-Ethyl-3-methylimidazolium hydroxide
MF C6 H11 N2 . H O
SR CA
LC STN Files: CA, CAPLUS, CASREACT, USPATFULL
CRN (65039-03-4)



● OH⁻

ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE
8 REFERENCES IN FILE CA (1907 TO DATE)
8 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 6 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN **244761-29-3** REGISTRY
ED Entered STN: 21 Oct 1999
CN Borate(1-), bis[ethanedioato(2-)-κO1,κO2]-, lithium, (T-4)-
(9CI) (CA INDEX NAME)
OTHER NAMES:
CN Lithium bis(oxalato)borate
CN Lithium bis(oxalato)borate(1-)
MF C4 B O8 . Li
CI CCS
SR CA
LC STN Files: CA, CAPLUS, CASREACT, CHEMCATS, CSCHEM, TOXCENTER, USPAT2, USPATFULL
CRN (125579-65-9)

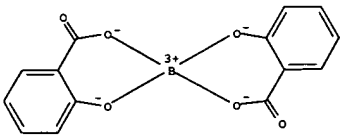


● Li⁺

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

141 REFERENCES IN FILE CA (1907 TO DATE)
2 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
143 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 7 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN **161589-07-7** REGISTRY
ED Entered STN: 17 Mar 1995
CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, lithium,
(T-4)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Borate(1-), bis[2-hydroxybenzoato(2-)-O1,O2]-, lithium, (T-4)-
OTHER NAMES:
CN AD 25
CN AD 25 (complex)
DR 496804-82-1
MF C14 H8 B O6 . Li
CI CCS
SR CA
LC STN Files: CA, CAPLUS, CASREACT, USPAT2, USPATFULL
CRN (38403-08-6)

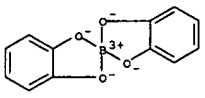


● Li⁺

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

22 REFERENCES IN FILE CA (1907 TO DATE)
22 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 8 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN **156762-86-6** REGISTRY
ED Entered STN: 03 Aug 1994
CN Borate(1-), bis[1,2-benzenediolato(2-)-κO,κO']-, lithium,
(T-4)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, lithium, (T-4)-
DR 244771-83-3
MF C12 H8 B O4 . Li
CI CCS, COM
SR CA
LC STN Files: CA, CAPLUS, CASREACT, DETHERM*, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
CRN (16986-25-7)

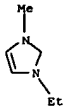


● Li⁺

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

26 REFERENCES IN FILE CA (1907 TO DATE)
26 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 9 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN **65039-09-0** REGISTRY
ED Entered STN: 16 Nov 1984
CN 1H-Imidazolium, 1-ethyl-3-methyl-, chloride (9CI) (CA INDEX NAME)
OTHER NAMES:
CN 1-Ethyl-3-methyl-1H-imidazolium chloride
CN 1-Ethyl-3-methylimidazolium chloride
CN 1-Methyl-3-ethylimidazolinium chloride
CN 1-Methyl-3-ethylimidazolium chloride
CN 3-Ethyl-1-methylimidazolium chloride
CN N-Methyl-N'-ethylimidazolium chloride
DR 140611-60-5
MF C6 H11 N2 . Cl
CI COM
LC STN Files: BEILSTEIN*, BIOSIS, CA, CAPLUS, CASREACT, CHEMCATS, CSCHEM, DETHERM*, GMELIN*, MSDS-OHS, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
CRN (65039-03-4)



● Cl⁻

ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

692 REFERENCES IN FILE CA (1907 TO DATE)
9 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
693 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 10 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN **10043-35-3** REGISTRY
ED Entered STN: 16 Nov 1984
CN Boric acid (H3BO3) (6CI, 8CI, 9CI) (CA INDEX NAME)
OTHER NAMES:
CN Basilit B
CN BC 140
CN Boracic acid
CN Boric acid
CN Boric acid (B(OH)3)
CN Borofax
CN Boron trihydroxide
CN Bortrac
CN CB BORid
CN Dia Flea-Mate
CN Dr.'s 1 Flea Terminator DF
CN Dr.'s 1 Flea Terminator DFPBO
CN Dr.'s 1 Flea Terminator DT
CN Dr.'s 1 Flea Terminator DTPBO
CN Flea Prufe
CN NSC 81726
CN Orthoboric acid
CN Orthoboric acid (B(OH)3)
CN Orthoboric acid (H3BO3)
CN Roach Away
CN Roach Prufe
CN Super Flea Eliminator
CN Trihydroxyborane
AR 11113-50-1
DR 12795-04-9
MF B H3 O3
CI COM
LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BIOSIS, BIOTECHNO, CA, CABA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DETHERM*, DIOGENES, DIPPR*, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MRCK*, MSDS-OHS, NIOSHTIC, PDLCOM*, PIRA, PROMT, PS, RTECS*, SPECINFO, TOXCENTER, TULSA, ULIDAT, USAN, USPAT2, USPATFULL,
VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

27542 REFERENCES IN FILE CA (1907 TO DATE)
1854 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
27591 REFERENCES IN FILE CAPLUS (1907 TO DATE)
1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

L3 ANSWER 11 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN **6153-56-6** REGISTRY
ED Entered STN: 16 Nov 1984
CN Ethanedioic acid, dihydrate (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Oxalic acid, dihydrate (8CI)
OTHER NAMES:
CN Oxalic acid dihydrate
MF C2 H2 O4 . 2 H2 O
LC STN Files: ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS, CA, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CSCHEM, DETHERM*, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MSDS-OHS, SPECINFO, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
CRN (144-62-7)

● 2 H_2O

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

315 REFERENCES IN FILE CA (1907 TO DATE)
2 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
315 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 12 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN **616-47-7** REGISTRY
ED Entered STN: 16 Nov 1984
CN 1H-Imidazole, 1-methyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Imidazole, 1-methyl- (6CI, 7CI, 8CI)
OTHER NAMES:
CN 1-Methyl-1H-imidazole
CN 1-Methylimidazole
CN Araldite DY 070
CN DY 070
CN N-Methylimidazole
CN N-methylimidazole
CN N1-Methylimidazole
CN NSC 88064
FS 3D CONCORD
DR 864745-22-2, 120418-32-8, 69723-05-3, 142504-34-5, 110069-11-9
MF C4 H6 N2
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU, EMBASE, GMELIN*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MSDS-OHS, NIOSHTIC, PIRA, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

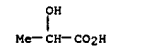
3634 REFERENCES IN FILE CA (1907 TO DATE)
219 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
3649 REFERENCES IN FILE CAPLUS (1907 TO DATE)
31 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

L3 ANSWER 13 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN 75-00-3 REGISTRY
ED Entered STN: 16 Nov 1984
CN Ethane, chloro- (8CI, 9CI) (CA INDEX NAME)
OTHER NAMES:
CN Aethylis
CN Aethylis chloridum
CN Anodynon
CN Chelen
CN Chlorene
CN Chlorethyl
CN Chloridum
CN Chloroethane
CN Chloryl
CN Chloryl Anesthetic
CN Chloryle anesthetic
CN Cloretilo
CN Dublofix
CN Ether chloratus
CN Ether hydrochloric
CN Ether muriatic
CN Ethyl chloride
CN F 160
CN Hydrochloric ether
CN Kelene
CN Monochlorethane
CN Monochloroethane
CN Muriatic ether
CN Narcotile
CN R 160
FS 3D CONCORD
MF C2 H5 Cl
CI COM
LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DIOGENES, DIPPR*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PDLCOM*, PROMT, PS, RTECS*, SCISEARCH, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, ULIDAT, USAN, USPAT2, USPATFULL, VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

4513 REFERENCES IN FILE CA (1907 TO DATE)
65 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
4515 REFERENCES IN FILE CAPLUS (1907 TO DATE)
5 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

L3 ANSWER 14 OF 14 REGISTRY COPYRIGHT 2006 ACS on STN
RN 50-21-5 REGISTRY
ED Entered STN: 16 Nov 1984
CN Propanoic acid, 2-hydroxy- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Lactic acid (7CI, 8CI)
OTHER NAMES:
CN (±)-Lactic acid
CN α-Hydroxypropanoic acid
CN α-Hydroxypropionic acid
CN 2-Hydroxy-2-methylacetic acid
CN 2-Hydroxypropanoic acid
CN 2-Hydroxypropionic acid
CN Biolac
CN Chem-Cast
CN DL-Lactic acid
CN dl-Lactic acid
CN E 270
CN Milk acid
CN NSC 367919
CN Purac FCC 80
CN Purac FCC 88
CN Tonsillosan
AR 849585-22-4
FS 3D CONCORD
DR 152-36-3, 598-82-3
MF C3 H6 O3
CI COM
LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CABA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DIOGENES, DIPPR*, DRUGU, EMBASE, ENCOMPLIT, ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IMSCOSEARCH, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, NIOSHTIC, PATDPASPC, PDLCOM*, PIRA, PROMT, PS, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, USAN, USPAT2, USPATFULL, VETU, VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

53297 REFERENCES IN FILE CA (1907 TO DATE)
1984 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
53385 REFERENCES IN FILE CAPLUS (1907 TO DATE)
1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

```
Connecting via Winsock to STN


Welcome to STN International!  Enter x:x

LOGINID:SSSPTA1600RXA

PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2


* * * * * * * * * *   Welcome to STN International   * * * * * * * * * *

 NEWS  1           Web Page URLs for STN Seminar Schedule - N. America
 NEWS  2           "Ask CAS" for self-help around the clock
 NEWS  3  DEC 05   CASREACT(R) - Over 10 million reactions available
 NEWS  4  DEC 14   2006 MeSH terms loaded in MEDLINE/LMEDLINE
 NEWS  5  DEC 14   2006 MeSH terms loaded for MEDLINE file segment of TOXCENTER
 NEWS  6  DEC 14   CA/CAplus to be enhanced with updated IPC codes
 NEWS  7  DEC 21   IPC search and display fields enhanced in CA/CAplus with the
                   IPC reform
 NEWS  8  DEC 23   New IPC8 SEARCH, DISPLAY, and SELECT fields in USPATFULL/
                   USPAT2
 NEWS  9  JAN 13   IPC 8 searching in IFIPAT, IFIUDB, and IFICDB
 NEWS 10  JAN 13   New IPC 8 SEARCH, DISPLAY, and SELECT enhancements added to
                   INPADOC
 NEWS 11  JAN 17   Pre-1988 INPI data added to MARPAT
 NEWS 12  JAN 17   IPC 8 in the WPI family of databases including WPIFV
 NEWS 13  JAN 30   Saved answer limit increased
 NEWS 14  JAN 31   Monthly current-awareness alert (SDI) frequency
                   added to TULSA
 NEWS 15  FEB 21   STN AnaVist, Version 1.1, lets you share your STN AnaVist
                   visualization results
 NEWS 16  FEB 22   Status of current WO (PCT) information on STN
 NEWS 17  FEB 22   The IPC thesaurus added to additional patent databases on STN
 NEWS 18  FEB 22   Updates in EPFULL; IPC 8 enhancements added
 NEWS 19  FEB 27   New STN AnaVist pricing effective March 1, 2006
 NEWS 20  FEB 28   MEDLINE/LMEDLINE reload improves functionality
 NEWS 21  FEB 28   TOXCENTER reloaded with enhancements
 NEWS 22  FEB 28   REGISTRY/ZREGISTRY enhanced with more experimental spectral
                   property data
 NEWS 23  MAR 01   INSPEC reloaded and enhanced
 NEWS 24  MAR 03   Updates in PATDPA; addition of IPC 8 data without attributes
 NEWS 25  MAR 08   X.25 communication option no longer available after June 2006

 NEWS EXPRESS  FEBRUARY 15 CURRENT VERSION FOR WINDOWS IS V8.01a,
               CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
               AND CURRENT DISCOVER FILE IS DATED 19 DECEMBER 2005.
               V8.0 AND V8.01 USERS CAN OBTAIN THE UPGRADE TO V8.01a AT
               http://download.cas.org/express/v8.0-Discover/

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 NEWS LOGIN    Welcome Banner and News Items
 NEWS PHONE    Direct Dial and Telecommunication Network Access to STN
 NEWS WWW      CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that
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  * * * * * * * * * * * * * * STN  Columbus  * * * * * * * * * * * * * * *

FILE 'HOME' ENTERED AT 09:13:16 ON 09 MAR 2006

=> fil reg
COST IN U.S. DOLLARS                                    SINCE FILE      TOTAL
                                                           ENTRY    SESSION
FULL ESTIMATED COST                                         0.21       0.21

FILE 'REGISTRY' ENTERED AT 09:13:25 ON 09 MAR 2006
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PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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STRUCTURE FILE UPDATES:    8 MAR 2006  HIGHEST RN 876273-86-8
DICTIONARY FILE UPDATES:   8 MAR 2006  HIGHEST RN 876273-86-8

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

  Please note that search-term pricing does apply when
  conducting SmartSELECT searches.

*********************************************************************
*                                                                   *
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added,   *
* effective March 20, 2005.  A new display format, IDERL, is now    *
* available and contains the CA role and document type information. *
*                                                                   *
*********************************************************************

Structure search iteration limits have been increased.  See HELP SLIMITS
for details.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document.  For information
on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end


=> screen 2040

L1    SCREEN CREATED

=>
Uploading C:\Program Files\Stnexp\Queries\QUERIES\09866926.str
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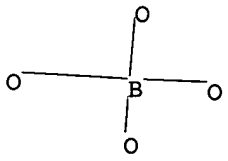
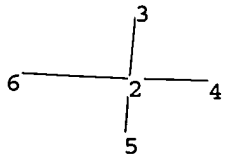




ring/chain nodes :
2 3 4 5 6
ring/chain bonds :
2-3 2-4 2-5 2-6
exact/norm bonds :
2-3 2-4 2-5 2-6

G1:C,H,X,Cb

Match level :
2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS

L2 STRUCTURE UPLOADED

=> que L2 AND L1

L3 QUE L2 AND L1

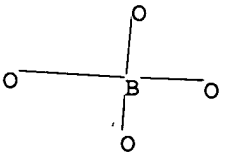
=> ....Testing the current file.... screen

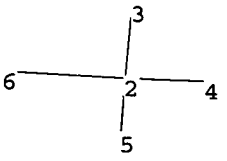
ENTER SCREEN EXPRESSION OR (END):end

=> screen 2040

L4 SCREEN CREATED

=>
Uploading C:\Program Files\Stnexp\Queries\QUERIES\09866926.str





ring/chain nodes :
2 3 4 5 6
ring/chain bonds :
2-3 2-4 2-5 2-6
exact/norm bonds :
2-3 2-4 2-5 2-6

G1:C,H,X,Cb

Match level :
2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS

L5 STRUCTURE UPLOADED

```
=> que L5 AND L4

L6   QUE L5 AND L4

=> s l6
SAMPLE SEARCH INITIATED 09:15:13 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -     1885 TO ITERATE

100.0% PROCESSED     1885 ITERATIONS                                  50 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:        35096 TO    40304
PROJECTED ANSWERS:            5566 TO     7754

L7           50 SEA SSS SAM L5 AND L4

=> s l6 full
FULL SEARCH INITIATED 09:15:18 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -    37720 TO ITERATE

100.0% PROCESSED    37720 ITERATIONS  (     1 INCOMPLETE)     6459 ANSWERS
SEARCH TIME: 00.00.01

L8         6459 SEA SSS FUL L5 AND L4

=> d his

     (FILE 'HOME' ENTERED AT 09:13:16 ON 09 MAR 2006)

     FILE 'REGISTRY' ENTERED AT 09:13:25 ON 09 MAR 2006
L1              SCREEN 2040
L2              STRUCTURE UPLOADED
L3              QUE L2 AND L1
L4              SCREEN 2040
L5              STRUCTURE UPLOADED
L6              QUE L5 AND L4
L7           50 S L6
L8         6459 S L6 FULL

=> s l3 subset=l8 full
FULL SUBSET SEARCH INITIATED 09:15:30 FILE 'REGISTRY'
FULL SUBSET SCREEN SEARCH COMPLETED -     6459 TO ITERATE

100.0% PROCESSED     6459 ITERATIONS                                 208 ANSWERS
SEARCH TIME: 00.00.01

L9          208 SEA SUB=L8 SSS FUL L2 AND L1

=> s l9 and caplus/lc
      49978644 CAPLUS/LC
L10        185 L9 AND CAPLUS/LC

=> s l9 not l10
L11         23 L9 NOT L10

=> d l11 1-23
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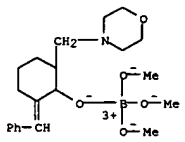
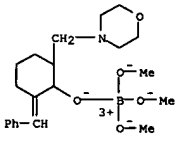
L11 ANSWER 1 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 852521-06-3 REGISTRY
ED Entered STN: 20 Jun 2005
CN INDEX NAME NOT YET ASSIGNED
MF C16 H74 B32 N16 O88 V10 Zn10
CI CCS, COM
SR CA

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

L11 ANSWER 2 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 795260-79-6 REGISTRY
ED Entered STN: 09 Dec 2004
CN Borate(1-), trimethoxy[rel-(1R,2R,6E)-2-(4-morpholinylmethyl)-6-(phenylmethylene)cyclohexanolato-κO]-, (T-4)- (9CI) (CA INDEX NAME)
MF C21 H33 B N O5
CI CCS, COM
SR CA



L11 ANSWER 3 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 795260-63-8 REGISTRY
ED Entered STN: 09 Dec 2004
CN Borate(1-), trimethoxy[rel-(1R,2S,6E)-2-(4-morpholinylmethyl)-6-(phenylmethylene)cyclohexanolato-κO]-, (T-4)- (9CI) (CA INDEX NAME)
MF C21 H33 B N O5
CI CCS, COM
SR CA



L11 ANSWER 4 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 792183-53-0 REGISTRY
ED Entered STN: 05 Dec 2004
CN Borate(1-), tetrakis(4-[2,2':5',2''-ter-1H-pyrrol]-5-ylphenolato-κO)-, (9CI) (CA INDEX NAME)
MF C72 H56 B N12 O4
CI CCS, COM
SR CA

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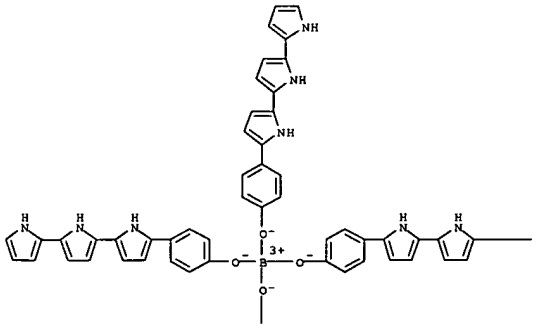

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L11 ANSWER 4 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN (Continued)

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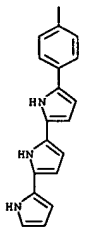
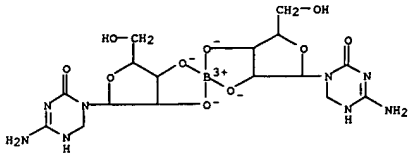


L11 ANSWER 5 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 792132-76-4 REGISTRY
ED Entered STN: 05 Dec 2004
CN Borate(1-), bis[4-amino-1-β-D-ribofuranosyl-1,3,5-triazin-2(1H)-onato(2-)-O2',O3']-, (T-4)- (9CI) (CA INDEX NAME)
MF C16 H24 B N8 O10
CI CCS, COM
SR CA



L11 ANSWER 6 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 790655-00-4 REGISTRY
ED Entered STN: 30 Nov 2004
CN Borate(1-), tetrakis(2-[2,2':5',2'':5'',2''':5''',2'''':4'''',2'''''-sexipyrazin]-5-ylbenzo[h]quinolin-10-olato-κO10)- (9CI) (CA INDEX NAME)
MF C148 H80 B N52 O4
CI CCS, COM
SR CA

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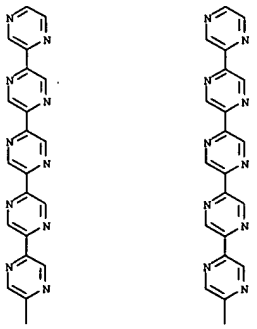

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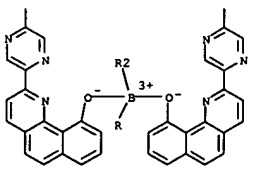

L11 ANSWER 6 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN (Continued)

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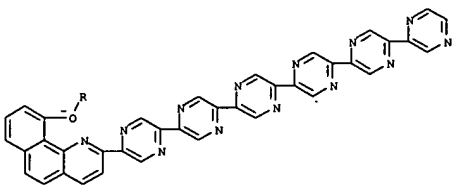

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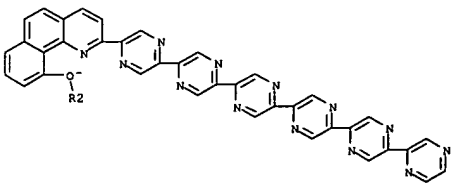

L11 ANSWER 7 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 788818-32-6 REGISTRY
ED Entered STN: 26 Nov 2004
CN Borate(1-),
bis[2-[2-(4-methyl[2,4':2',4'':2'',4''':2''',4'''':2'''',4''''
':2'''''',4''''':2''''''',4'''''':2'''''',4''''''':2'''''''-novi-1H-imidazol]-
2'''''''-yl)-2-phenylethenyl]-8-quinolinolato-κO8]bis(8-
quinolinolato-κO8)-, (T-4)- (9CI) (CA INDEX NAME)
MF C108 H76 B N40 O4
CI CCS, COM
SR CA

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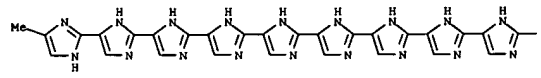

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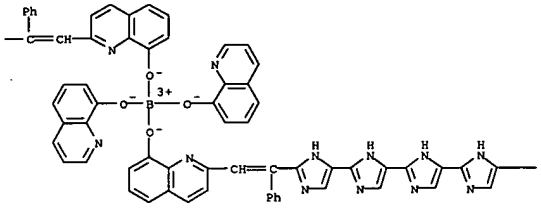

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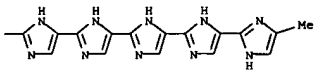

L11 ANSWER 7 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN (Continued)

L11 ANSWER 8 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 780750-08-5 REGISTRY
ED Entered STN: 15 Nov 2004
CN Borate(1-),
tetrakis[4-(1,1'-dimethyl-3,3',4,4',5'-pentaphenyl[2,2'-bi-1H-
pyrrol]-5-yl)-8-quinolinolato-κO8]- (9CI) (CA INDEX NAME)
MF C196 H144 B N12 O4
CI CCS, COM
SR CA

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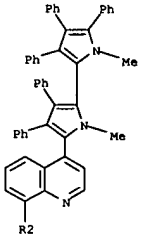

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

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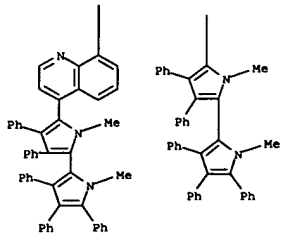

L11 ANSWER 9 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 778571-66-7 REGISTRY
ED Entered STN: 11 Nov 2004
CN Borate(1-),
tetrakis[2-[2-[5-(4-nitrophenyl)[2,5':2',5'':2'',5''':2''',5''
''-quinquethiazol]-2''''-yl]ethenyl]-8-quinolinolato-κO8]- (9CI)
(CA INDEX NAME)
MF C128 H64 B N28 O12 S20
CI CCS, COM
SR CA

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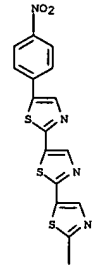

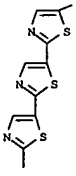

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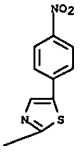

L11 ANSWER 9 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN (Continued)

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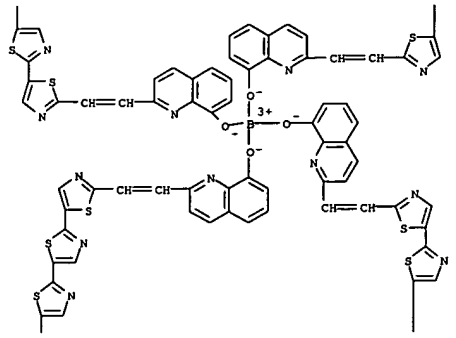

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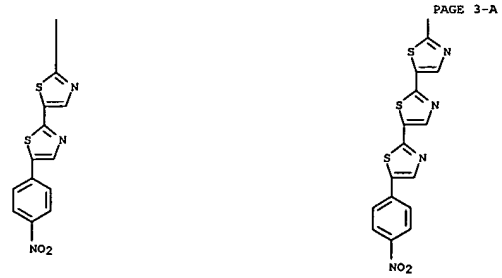
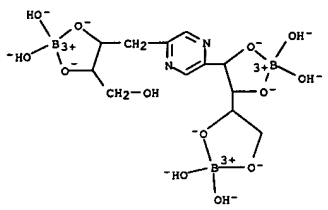


L11 ANSWER 10 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 777846-21-6 REGISTRY
ED Entered STN: 10 Nov 2004
CN Borate(3-), hexahydroxy[μ3-{(1R,2S,3R)-1-{5-{(2S,3R)-2,3-di(hydroxy-κO)-4-hydroxybutyl]pyrazinyl}-1,2,3,4-butanetetrolato(6-)-κO1,κO2:κO3,κO4]]tri- (9CI) (CA INDEX NAME)
MF C12 H20 B3 N2 O13
CI CCS, COM
SR CA



L11 ANSWER 11 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 772335-27-0 REGISTRY
ED Entered STN: 31 Oct 2004
CN Borate(1-), tetrakis[2-[2-[5-(4-methoxyphenyl)[2,5':2',5''-teroxazol]-2''-yl]ethenyl]phenolato-κO]- (9CI) (CA INDEX NAME)
MF C96 H64 B N12 O20
CI CCS, COM
SR CA

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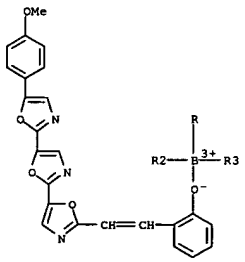

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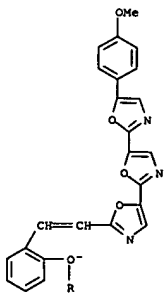

L11 ANSWER 11 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN (Continued)

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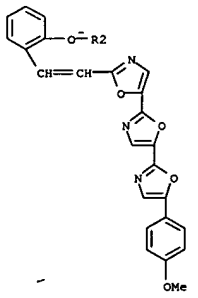

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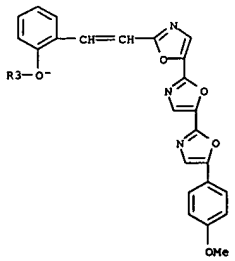

L11 ANSWER 12 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 749844-94-8 REGISTRY
ED Entered STN: 23 Sep 2004
CN Borate(1-),
tetrakis[4-(1,1'-dimethyl-3,3',4,4',5'-pentaphenyl[2,2'-bi-1H-pyrrol]-5-yl)benzo[h]quinolin-10-olato-κO10]- (9CI) (CA INDEX NAME)
MF C212 H152 B N12 O4
CI CCS, COM
SR CA

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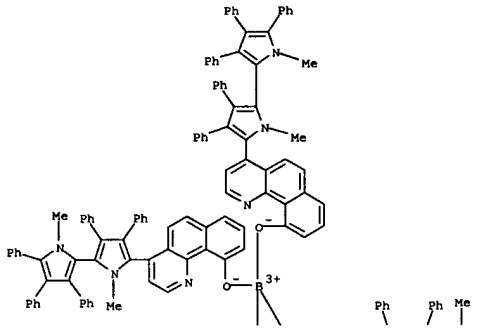

L11 ANSWER 12 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN (Continued)

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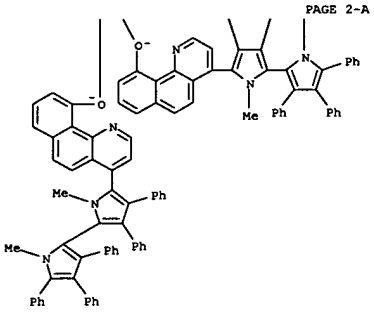

L11 ANSWER 13 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 741667-63-0 REGISTRY
ED Entered STN: 09 Sep 2004
CN Borate(1-), tetrakis(2-[2,3'-bipyridin]-5'-yl-8-quinolinolato-κO8)-(9CI) (CA INDEX NAME)
MF C76 H48 B N12 O4
CI CCS, COM
SR CA

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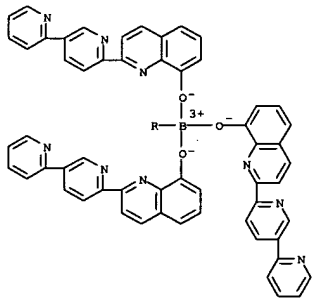

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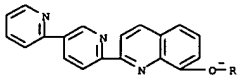

L11 ANSWER 14 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 741246-32-2 REGISTRY
ED Entered STN: 08 Sep 2004
CN Borate(1-),
tetrakis[7-[2-[5-(4-nitrophenyl)[2,5':2',5'':2'',5''':2''',5''''-quinquethiazol]-2''''-yl]ethenyl]benzo[f]quinolin-5-olato-κO5]-(9CI) (CA INDEX NAME)
MF C144 H72 B N28 O12 S20
CI CCS, COM
SR CA

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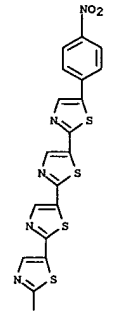

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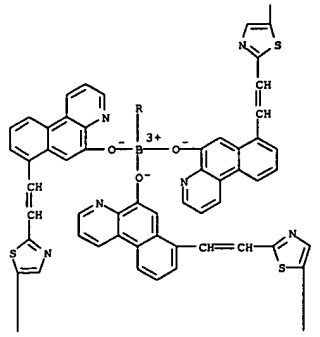

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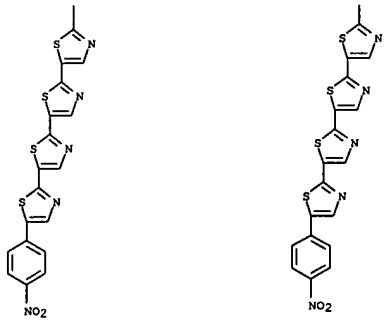

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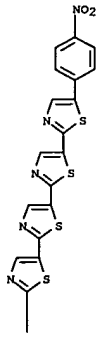

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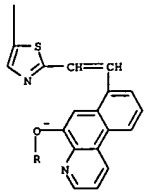

L11 ANSWER 15 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 737749-88-1 REGISTRY
ED Entered STN: 02 Sep 2004
CN Borate(2-), [μ-[1-(hydroxy-κO)piperidinato-κN]][1-(hydroxy-κO)piperidine][μ-[orthoborato(3-)-κO:κO']]di-μ-oxophenyl[(phenylboronic acid-κO) bimol. monoanhydridato(2-)]tri- (9CI) (CA INDEX NAME)
MF C28 H36 B6 N2 O10
CI CCS, COM
SR CA

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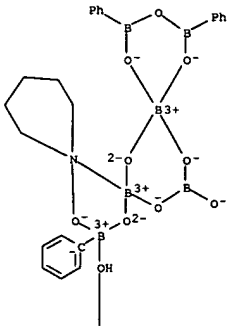

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L11 ANSWER 16 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 733735-50-7 REGISTRY
ED Entered STN: 27 Aug 2004
CN Borate(1-), tetrakis[5-[2-[5'-(4-pyridinyl)[2,2'-bi-1,3,4-oxadiazol]-5-yl]ethenyl]-8-quinolinolato-κO8]- (9CI) (CA INDEX NAME)
MF C80 H44 B N24 O12
CI CCS, COM
SR CA

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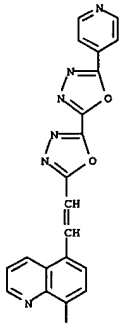

L11 ANSWER 16 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN (Continued)

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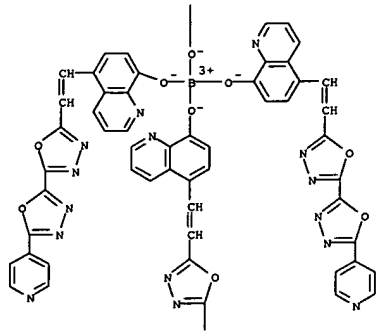

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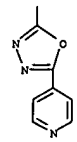
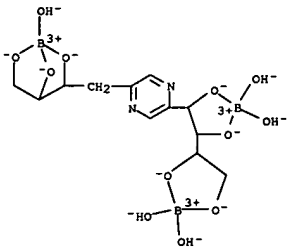


L11 ANSWER 17 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 689215-18-7 REGISTRY
ED Entered STN: 03 Jun 2004
CN Borate(3-), pentahydroxy[μ3-[(1R,2S,3R)-1-[5-[(2S,3R)-2,3,4-tri(hydroxy-κO)butyl]pyrazinyl]-1,2,3,4-butanetetrolato(7-)-κO1,κO2:κO3,κO4]]tri- (9CI) (CA INDEX NAME)
MF C12 H18 B3 N2 O12
CI CCS, COM
SR CA

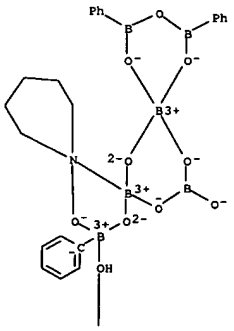


L11 ANSWER 18 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 433330-51-9 REGISTRY
ED Entered STN: 25 Jun 2002
CN Borate(2-), [μ-[1-(hydroxy-κO)piperidinato-κN]][1-(hydroxy-κO)piperidine][μ-[orthoborato(3-)-κO:κO']]di-μ-oxophenyl[(phenylboronic acid-κO) bimol. monoanhydridato(2-)]tri-, dihydrogen, compd. with 1-hydroxypiperidine (1:1) (9CI) (CA INDEX NAME)
MF C28 H36 B6 N2 O10 . C5 H11 N O . 2 H
CI COM
SR CA

CM 1

CRN 433330-50-8 (737749-88-1)
CMF C28 H36 B6 N2 O10 . 2 H
CCI CCS

PAGE 1-A



PAGE 2-A

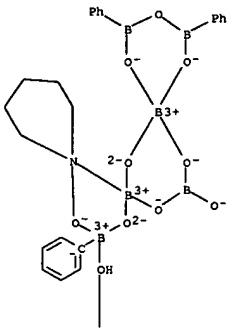
●2 H+

L11 ANSWER 18 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN (Continued)
CM 2

CRN 4801-58-5
CMF C5 H11 N O

L11 ANSWER 19 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 433330-50-8 REGISTRY
ED Entered STN: 25 Jun 2002
CN Borate(2-), [μ-[1-(hydroxy-κO)piperidinato-κN]][1-(hydroxy-κO)piperidine][μ-[orthoborato(3-)-κO:κO']]di-μ-oxophenyl[(phenylboronic acid-κO) bimol. monoanhydridato(2-)]tri-, dihydrogen (9CI) (CA INDEX NAME)
MF C28 H36 B6 N2 O10 . 2 H
CI CCS, COM
SR CA
CRN (737749-88-1)

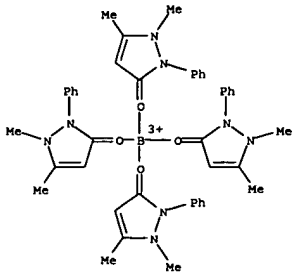
PAGE 1-A



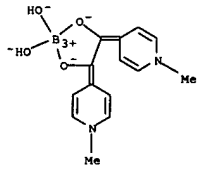
PAGE 2-A

●2 H+

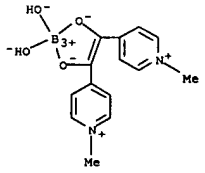
L11 ANSWER 20 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 84663-12-7 REGISTRY
ED Entered STN: 16 Nov 1984
CN Boron(3+), tetrakis(1,2-dihydro-1,5-dimethyl-2-phenyl-3H-pyrazol-3-one-O)-, (T-4)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 3H-Pyrazol-3-one, 1,2-dihydro-1,5-dimethyl-2-phenyl-, boron complex
MF C44 H48 B N8 O4
CI CCS, COM



L11 ANSWER 21 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 51095-90-0 REGISTRY
ED Entered STN: 16 Nov 1984
CN Borate(1-), [1,2-bis(1-methyl-4(1H)-pyridinylidene)-1,2-ethanediolato(2-)-O,O']dihydroxy-, (T-4)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 1,2-Ethanediol, 1,2-bis(1-methyl-4(1H)-pyridinylidene)-, boron complex
MF C14 H16 B N2 O4
CI CCS, COM



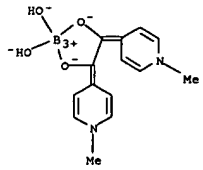
L11 ANSWER 22 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 51095-89-7 REGISTRY
ED Entered STN: 16 Nov 1984
CN Boron(1+), [[4,4'-(1,2-dihydroxy-1,2-ethenediyl)bis[1-methylpyridiniumato]](2-)-O,O']dihydroxy-, (T-4)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Pyridinium, 4,4'-(1,2-dihydroxy-1,2-ethenediyl)bis[1-methyl-, boron complex
MF C14 H16 B N2 O4
CI CCS, COM



L11 ANSWER 23 OF 23 REGISTRY COPYRIGHT 2006 ACS on STN
RN 6985-01-9 REGISTRY
ED Entered STN: 16 Nov 1984
CN Boron(1+), [[4,4'-(1,2-dihydroxy-1,2-ethenediyl)bis[1-methylpyridiniumato]](2-)-O,O']dihydroxy-, (T-4),
(T-4)-[1,2-bis(1-methyl-4(1H)-pyridinylidene)-1,2-ethanediolato(2-)-O,O']dihydroxyborate(1-)
(9CI)
(CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 1,2-Ethanediol, 1,2-bis(1-methyl-4(1H)-pyridinylidene)-, boron complex
CN Pyridinium, 4,4'-(1,2-dihydroxy-1,2-ethenediyl)bis[1-methyl-, boron complex
MF C14 H16 B N2 O4 . C14 H16 B N2 O4

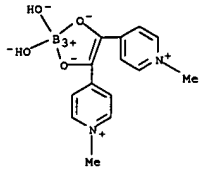
CM 1

CRN 51095-90-0
CMF C14 H16 B N2 O4
CCI CCS



CM 2

CRN 51095-89-7
CMF C14 H16 B N2 O4
CCI CCS



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                                                           ENTRY    SESSION
FULL ESTIMATED COST                                       256.56     256.77

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L2               STRUCTURE UPLOADED
L3               QUE L2 AND L1
L4               SCREEN 2040
L5               STRUCTURE UPLOADED
L6               QUE L5 AND L4
L7            50 S L6
L8          6459 S L6 FULL
L9           208 S L3 FULL SUB=L8
L10          185 S L9 AND CAPLUS/LC
L11           23 S L9 NOT L10

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L12         105 L10

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L12 ANSWER 1 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:980559 CAPLUS
DOCUMENT NUMBER: 144:12416
TITLE: Thermochemistry of triimidazolium nonaborate
AUTHOR(S): Liu, Zhi-Hong; Zhang, Wen-Juan
CORPORATE SOURCE: School of Chemistry and Materials Science, Shaanxi Normal University, Xi'an, 710062, Peop. Rep. China
SOURCE: Thermochimica Acta (2005), 436(1-2), 156-158
CODEN: THACAS; ISSN: 0040-6031
PUBLISHER: Elsevier B.V.
DOCUMENT TYPE: Journal
LANGUAGE: English

AB The crystalline triimidazolium nonaborate [C3H5N2]3[B9O12(OH)6] has been prepared and identified by XRD, TG, elemental anal., and chemical anal. The molar enthalpy of solution of [C3H5N2]3[B9O12(OH)6] in 0.9996 mol dm-3 HCl(aq) and of imidazole in (HCl + H3BO3) (aq) were determined With the incorporation of the enthalpies of solution of H3BO3 in HCl(aq), and the standard molar enthalpies of formation of C3H4N2(s), H3BO3(s), and H2O(l), the standard molar enthalpy of formation of -7116 ± 7 kJ mol-1 of [C3H5N2]3[B9O12(OH)6] was obtained.

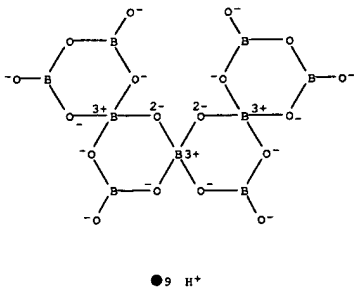
IT **273750-78-0**
RL: PRP (Properties); RCT (Reactant); RACT (Reactant or reagent)
(enthalpy of formation of triimidazolium nonaborate from solution calorimetry)

RN 273750-78-0 CAPLUS
CN Borate(9-), bis[μ-[orthoborato(3-)-κO:κO']]di-μ-oxobis[μ-oxotetraoxodiborato(4-)]tri-, nonahydrogen, compd. with 1H-imidazole (1:9) (9CI) (CA INDEX NAME)

CM 1

CRN 273750-76-8
CMF B9 O18 . 9 H
CCI CCS

L12 ANSWER 1 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



●9 H+

CM 2

CRN 288-32-4
CMF C3 H4 N2

REFERENCE COUNT: 10 THERE ARE 10 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 2 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:491174 CAPLUS
DOCUMENT NUMBER: 144:157759
TITLE: Tailor-made ionic liquids
AUTHOR(S): Jork, C.; Kristen, C.; Pieraccini, D.; Stark, A.; Chiappe, C.; Beste, Y. A.; Arlt, W.
CORPORATE SOURCE: Institut fuer Verfahrenstechnik, Fachgebiet Thermodynamik und Thermische Verfahrenstechnik, Technische Universitaet Berlin, Berlin, 10623, Germany
SOURCE: Journal of Chemical Thermodynamics (2005), 37(6), 537-558
CODEN: JCTDAF; ISSN: 0021-9614
PUBLISHER: Elsevier Ltd.
DOCUMENT TYPE: Journal
LANGUAGE: English

AB This article presents a first consequent thermodn. optimization of ionic liqs. (IL) as entrainers in the separation upon distillation of both an azeotropic aqueous (THF + water) and a close-boiling aromatic test system (methylcyclohexane + toluene) on the basis of COSMO-RS predictions. The use of this method allows for the preselection from the large pool of available IL. Thus, favorable structural variations were identified and used for tailoring IL entrainers. For the prediction of activity coeffs. with COSMO-RS, the use of different conformations of the components, derived from conformational analyses, leads to varying results. The simulations showed that the influence of conformations of the volatile components and the ionic liqs. depends largely on the type of the phase equilibrium, which is investigated.
The approach to tailor ionic liqs. as additives for separation science starts with the prediction of the activity coeffs. at infinite dilution The simulation indicated that a higher degree of branching or longer alkyl substituents on the cation, as well as a low nucleophilicity of the anion decreases both selectivity and capacity in the polar test mixture However, COSMO-RS calcns. for the non-polar mixture showed that the selection of an entrainer for this system is more complicated, because - contrarily to (THF + water) - structural variations of the IL entrainer cause converse changes in selectivity and capacity: while the selectivity for toluene increases with a lower degree of branching and a shorter alkyl substituent of the cation as well as with a lower nucleophilicity of the anion, these properties decrease the capacity. In this work, the most favorable IL entrainers were synthesized and the separation factors of the test systems were exptl. validated at finite dilution

IT **566135-35-1P**
RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)
(phase equilibrium and constituent activity in water/organic solvent/ionic liquid binary and ternary mixts.)

RN 566135-35-1 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[ethanedioato(2-)-κO1,κO2]borate(1-) (9CI) (CA INDEX NAME)

L12 ANSWER 2 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

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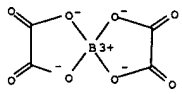
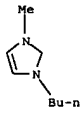
CRN 125579-65-9
CMF C4 B O8
CCI CCS



CM 2

CRN 80432-08-2
CMF C8 H15 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE
REFERENCE COUNT: 29 THERE ARE 29 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 3 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:351310 CAPLUS
DOCUMENT NUMBER: 143:317708
TITLE: Thermal decomposition mechanism and non-isothermal kinetics of the polyoxometalate of ciprofloxacin with 12-tungstoboric acid
AUTHOR(S): Wang, Dunjia; Fang, Zhengdong; Han, Deyan
CORPORATE SOURCE: Department of Chemistry and Environmental Engineering, Hubei Normal University, Huangshi, 435002, Peop. Rep. China
SOURCE: Rare Metals (Beijing, China) (2005), 24(1), 15-21
CODEN: RARME8; ISSN: 1001-0521
PUBLISHER: Rare Metals
DOCUMENT TYPE: Journal
LANGUAGE: English

AB The polyoxometalate complex (CPFX·HCl)4H5BW12O40·12H2O was prepared in aqueous solution for the 1st time, and characterized by elemental anal., IR spectrum, and TG-DTG. The TG-DTG curves showed that its thermal decomposition was a four-step process consisting of the simultaneous collapse of Keggin anion. The intermediate and residue of the decomposition were identified by mean of TG-DTG, IR, and XRD technique. The nonisothermal kinetic data were analyzed by the Achar method and Coats-Redfern method. The apparent activation energy (E) and the pre-exponential factor (ln A) of each decomposition were obtained. The most probable thermal decomposition reaction mechanisms are proposed by comparison of the kinetic parameters. The kinetic equation for both the 2nd stage and the 3rd stage can be expressed as dα/dt = Ae-E/RT·(1 - α)2, and the 4th stage dα/dt = Ae-E/RT·(1 - α). And their mathematic expressions of the kinetic compensation effects of thermal decomposition reaction were also determined

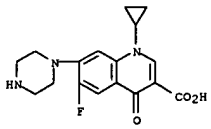
IT **864948-78-7P**
RL: CPS (Chemical process); PEP (Physical, engineering or chemical process); PRP (Properties); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)
(preparation and nonisothermal kinetics of thermal decomposition of)

RN 864948-78-7 CAPLUS
CN Tungstate(5-), tetracosa-μ-oxododecaoxo[μ12-[tetrahydroxyborato(5-)-κO:κO:κO:κO':κO':κO'':.kappa.O'':κO'':κO''':κO''':κO''']]dodeca-, pentahydrogen, compd. with 1-cyclopropyl-6-fluoro-1,4-dihydro-4-oxo-7-(1-piperazinyl)-3-quinolinecarboxylic acid monohydrochloride (1:4), dodecahydrate (9CI) (CA INDEX NAME)

CM 1

CRN 93107-08-5
CMF C17 H18 F N3 O3 . Cl H

L12 ANSWER 3 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

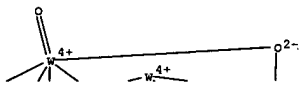


● HCl

CM 2

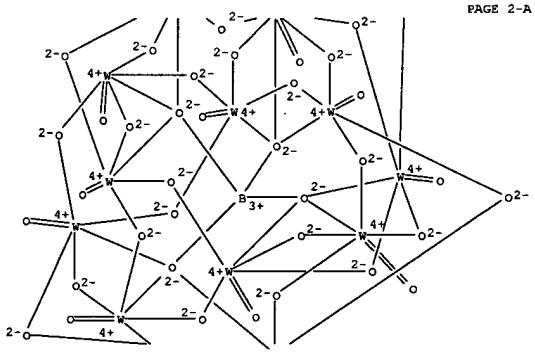
CRN 12297-12-0
CMF B 040 W12 . 5 H
CCI CCS

PAGE 1-A

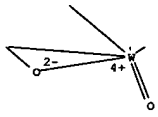


L12 ANSWER 3 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

PAGE 2-A



PAGE 3-A



●5 H^+

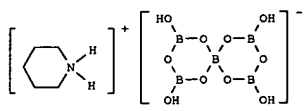
REFERENCE COUNT: 18 THERE ARE 18 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 4 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:297601 CAPLUS
DOCUMENT NUMBER: 142:356086
TITLE: Piperidine condensed borate salt, its compositions, epoxy resin hardeners thereof, and epoxy resin compositions containing them
INVENTOR(S): Obayashi, Akira; Haraguchi, Kazutoshi; Maki, Hiroshi
PATENT ASSIGNEE(S): Dainippon Ink and Chemicals, Inc., Japan; Kawamura Institute of Chemical Research
SOURCE: Jpn. Kokai Tokkyo Koho, 11 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2005089317	A2	20050407	JP 2003-321197	20030912
PRIORITY APPLN. INFO.:			JP 2003-321197	20030912

GI



I

AB The title salt I, giving heat-resistant epoxy resins as a hardener, is prepared Thus, 30 g H3BO3 was treated with 8.3 g piperidine in DMF, washed, and dried to give 21.1 g I. A composition of Epiclon 850 (epoxy resin) 100, I 16, MeOH 80, and MEK 80 g was heated to 50° for 3 h, applied on an Al foil, dried, and heated to give a cured product showing light transmittance 92% and Tg 285°.

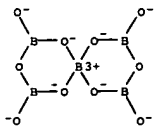
IT **12548-84-4P**
RL: IMF (Industrial manufacture); MOA (Modifier or additive use); PRP (Properties); RCT (Reactant); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(piperidine condensed borate salt as epoxy resin hardener for good heat resistance)

RN 12548-84-4 CAPLUS
CN Borate(5-), bis[μ-oxotetraoxodiborato(4-)]-, (T-4)-, pentahydrogen, compd. with piperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 92258-67-8
CMF B5 O10 . 5 H
CCI CCS

L12 ANSWER 4 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



●5 H^+

CM 2

CRN 110-89-4
CMF C5 H11 N

L12 ANSWER 5 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2005:267909 CAPLUS
DOCUMENT NUMBER: 143:18668
TITLE: Synthesis of elliptical vanadoborates housing bimetallic centers [Zn4(B2O4H2)(V10B28O74H8)]8- and [Mn4(C2O4)(V10B28O74H8)]10-
AUTHOR(S): Wu, Mingmei; Law, Teresa S-C.; Sung, Herman H-Y.; Cai, Jiwen; Williams, Ian D.
CORPORATE SOURCE: Department of Chemistry, Hong Kong University of Science and Technology, Hong Kong, Peop. Rep. China
SOURCE: Chemical Communications (Cambridge, United Kingdom) (2005), (14), 1827-1829
CODEN: CHCOFS; ISSN: 1359-7345
PUBLISHER: Royal Society of Chemistry
DOCUMENT TYPE: Journal
LANGUAGE: English
OTHER SOURCE(S): CASREACT 143:18668
AB The hydrothermal preparation and crystal structures of 3 new vanadoborate compds. with elliptical (V10B28O74H8) clusters, [Zn(en)2]2 [Zn(OH2)(en)]4 [Zn4(B2O4H2)(BO3H)2(V10B28O74H8)]·10H2O (1), K6(H3O)8[Mn4(C2O4)(V10B28O74H8)(B4O9H2)]·24H2O (2), and (H3O)14[Mn4(C2O4)3(V10B28O74H8)]·32H2O (3) is described. The clusters contain pairs of bimetallic Zn2 or Mn2 units.
IT **852461-51-9P**
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (preparation and crystal structure of 2-dimensional polymeric)
RN 852461-51-9 CAPLUS
CN Vanadate(10-), (di-μ-hydroxydiborate)bis[μ14-[hexakis[μ-[orthoborato(3-)-κO:κO']]hexa-μ-oxooctaoxooctaborato(22-)]]di-μ-oxodi-μ3-oxo[tetraaquaoctakis(1,2-ethanediamine-κN,κN')bis[μ-[orthoborato(3-)-κO:κO']]decazincate]deca-, decahydrogen, decahydrate (9CI) (CA INDEX NAME)

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
REFERENCE COUNT: 19 THERE ARE 19 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 6 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2004:841950 CAPLUS
DOCUMENT NUMBER: 141:358096
TITLE: Ink-jet recording media with high ink absorbability and providing high-resolution water-resistant images
INVENTOR(S): Ito, Kengo; Uehara, Masaharu
PATENT ASSIGNEE(S): Sony Corp., Japan; Boron International K. K.
SOURCE: Jpn. Kokai Tokkyo Koho, 12 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2004284085	A2	20041014	JP 2003-76600	20030319
PRIORITY APPLN. INFO.:			JP 2003-76600	20030319

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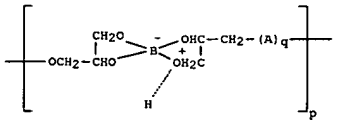


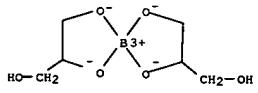
I

AB In the media, ink-receiving layers formed on substrates contain binders capable of forming chemical bonds with the boron compds. and fillers, the surface of which are modified with semipolar organic boron compds. Preferably, the boron compds. are expressed by I [A = Xl-Ym-Zn; X, Z = C≤100 oxygen-containing hydrocarbyl bearing ether terminal group; Y = OCORCO; R = C1-34 hydrocarbyl; l, m, n = 0, 1; P = 10-1000; q ≥1] or their derivs.
IT **773868-49-8**
RL: TEM (Technical or engineered material use); USES (Uses) (coating on filler; ink-jet recording medium containing organoboron-surface-modified filler and binder forming chemical bond with organoboron)
RN 773868-49-8 CAPLUS
CN Borate(1-), bis[1,2,3-propanetriolato(2-)-κO1,κO2]-, hydrogen, (T-4)-, polymer with 1-ethenyl-2-pyrrolidinone (9CI) (CA INDEX NAME)

CM 1

CRN 49625-59-4
CMF C6 H12 B O6 . H
CCI CCS

L12 ANSWER 6 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



● H^+

CM 2

CRN 88-12-0
CMF C6 H9 N O

L12 ANSWER 7 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:534653 CAPLUS
DOCUMENT NUMBER: 142:47884
TITLE: Synthesis and anti-liver cancer activity of 5-fluorouracil salt of 12-tungstoboric acid
AUTHOR(S): Li, Juan; Li, Jing; Qi, Yan-fei; Wang, Hong-Fang; Wang, En-Bo; Hu, Chang-Wen; Xu, Lin; Wu, Xin-Yu
CORPORATE SOURCE: Inst. Polyoxometalate Chem., Fac. Chem., Northeast Normal Univ., Changchun, 130021, Peop. Rep. China
SOURCE: Gaodeng Xuexiao Huaxue Xuebao (2004), 25(6), 1010-1012
CODEN: KTHPDM; ISSN: 0251-0790
PUBLISHER: Gaodeng Jiaoyu Chubanshe
DOCUMENT TYPE: Journal
LANGUAGE: Chinese
OTHER SOURCE(S): CASREACT 142:47884

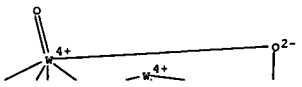
AB 5-Fluorouracil salt (WBF) of 12-tungstoboric acid with Keggin structure was synthesized and its structure was characterized by IR, 1H NMR, 183W NMR and elementary anal. The MTT exptl. results show that WBF salt has a higher anti-tumor activity for liver cancer cells in vitro and lower toxicity than the unsalted 5-fluorouracil. The acute toxicity of WBF was measured. The half toxicity dose LD50 of WBF and 5-fluorouracil orally administrated to mice is 1117.4 and 220 mg/kg, resp.

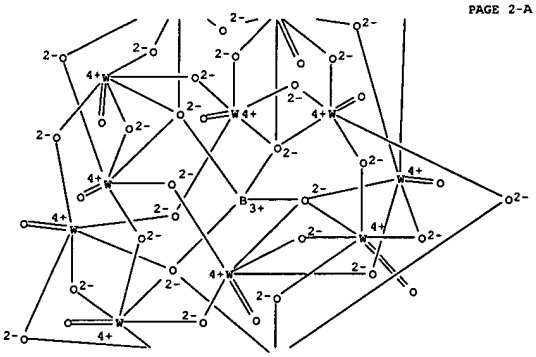
IT **803738-68-3P**
RL: BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation and anti-liver cancer activity of 5-fluorouracil salt of 12-tungstoboric acid)

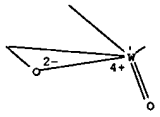
RN 803738-68-3 CAPLUS
CN Tungstate(5-), tetracosa-μ-oxododecaoxo[μ12-[tetrahydroxyborato(5-)-κO:κO:κO:κO':κO':κO':κO'':.kappa.O'':κO'':κO''':κO''':κO''']]dodeca-, pentahydrogen, compd. with 5-fluoro-2,4(1H,3H)-pyrimidinedione (1:5) (9CI)
(CA INDEX NAME)

CM 1

CRN 12297-12-0
CMF B O40 W12 . 5 H
CCI CCS







●5 H+

CM 2

CRN 51-21-8
CMF C4 H3 F N2 O2

L12 ANSWER 8 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2004:525791 CAPLUS
DOCUMENT NUMBER: 141:286190
TITLE: 4-Methylpyridinium bis(pyrocatecholato-κ2O,O')borate catechol solvate
AUTHOR(S): Baber, R. Angharad; Charmant, Jonathan P. H.; Moore, Jonathan D.; Norman, Nicholas C.; Orpen, A. Guy
CORPORATE SOURCE: School of Chemistry, University of Bristol, Bristol, BS8 1TS, UK
SOURCE: Acta Crystallographica, Section E: Structure Reports Online (2004), E60(7), o1140-o1142
CODEN: ACSEBH; ISSN: 1600-5368
URL: http://journals.iucr.org/e/journalhomepage.html
PUBLISHER: International Union of Crystallography
DOCUMENT TYPE: Journal; (online computer file)
LANGUAGE: English

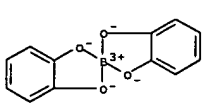
AB Unlike the previously reported salts of the 4-methylpyridinium cation and the bis(pyrocatecholato)borate anion [Clegg et al. (1998). Acta Crystalline C54, 1875-1880], the title compound, C6H8N+·C12H8BO4-·C6H6O2, is a solvate containing a mol. of catechol. The crystal packing is influenced by N-H···O and O-H···O H bonds. Crystallog. data are given.

IT **758722-93-9**
RL: PRP (Properties)
(crystal structure of)

RN 758722-93-9 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with 1,2-benzenediol and 4-methylpyridine (1:1:1) (9CI)
(CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H+

CM 2

CRN 120-80-9
CMF C6 H6 O2

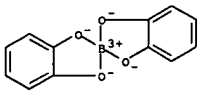
CM 3

CRN 108-89-4
CMF C6 H7 N

IT **219702-48-4**
RL: PRP (Properties)
(preparation and 11B NMR of)
RN 219702-48-4 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with 4-methylpyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H^+

CM 2

CRN 108-89-4
CMF C6 H7 N

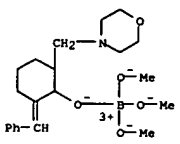
REFERENCE COUNT: 8 THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE

FORMAT

ACCESSION NUMBER: 2004:401384 CAPLUS
DOCUMENT NUMBER: 141:206835
TITLE: Diastereoselective reduction of cyclic bioactive Mannich ketones
AUTHOR(S): Lorandd, Tamas; Osz, Erzsebet; Kispal, Gyula; Nagy, Gergely; Weckert, Edgar; Luebbert, Daniel; Meents, Alke; Kocsis, Bela; Prokai, Laszlo
CORPORATE SOURCE: Faculty of Medicine, Department of Biochemistry and Medical Chemistry, University Pecs, Pecs, H-7601, Hung.
SOURCE: ARKIVOC (Gainesville, FL, United States) (2004), (7), 34-52
CODEN: AGFUAR
URL: http://www.arkat-usa.org/ark/journal/2004/Antus/S A-837B/837B.pdf
PUBLISHER: Arkat USA Inc.
DOCUMENT TYPE: Journal; (online computer file)
LANGUAGE: English
OTHER SOURCE(S): CASREACT 141:206835

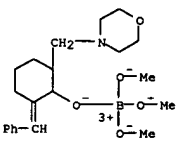
AB The reduction of cyclic Mannich ketones having antibacterial activity showed poor stereoselectivity with L-Selectride, sodium trimethoxyborohydride and diisobutylaluminum hydride, while lithium aluminum hydride and, in particular, sodium borohydride often yielded a single stereoisomer. The size of the ring strongly influenced the stereocomposition of the reaction mixts. An increased preference for the trans isomer was attributed to a weak intramol. hydrogen bond between the OH and the N, as demonstrated by X-ray crystallog. After reduction, the antibacterial activity of the product decreased dramatically ("S-form") but was not completely abolished in deep rough mutants of Gram-neg. bacteria.

IT **741725-62-2 742077-99-2**
RL: PRP (Properties)
(calculated heat of formation of intermediates from diastereoselective reduction of cyclic Mannich ketones)
RN 741725-62-2 CAPLUS
CN Borate(1-), trimethoxy[rel-(1R,2S,6E)-2-(4-morpholinylmethyl)-6-(phenylmethylene)cyclohexanolato-κO]-, sodium, (T-4)- (9CI) (CA INDEX NAME)



● Na^+

RN 742077-99-2 CAPLUS
CN Borate(1-), trimethoxy[rel-(1R,2R,6E)-2-(4-morpholinylmethyl)-6-(phenylmethylene)cyclohexanolato-κO]-, sodium, (T-4)- (9CI) (CA INDEX NAME)

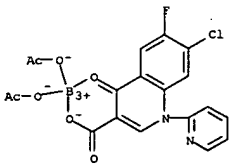


● Na^+

REFERENCE COUNT: 16 THERE ARE 16 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE

FORMAT

L12 ANSWER 10 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2004:253357 CAPLUS
DOCUMENT NUMBER: 141:106352
TITLE: Synthesis and antibacterial activity of N-pyridine quinolone derivative
AUTHOR(S): Wang, Dun-jia; Huang, Ling
CORPORATE SOURCE: Department of Chemistry and Environmental Engineering, Hubei Normal University, Huangshi, 435002, Peop. Rep. China
SOURCE: Huaxue Shiji (2004), 26(1), 47-49
CODEN: HUSHDR; ISSN: 0258-3283
PUBLISHER: Huagongbu Huaxue Shiji Xinsizhan
DOCUMENT TYPE: Journal
LANGUAGE: Chinese
OTHER SOURCE(S): CASREACT 141:106352
AB 1-(2-Pyridyl)-7-chloro-6-fluoro-1,4-dihydro-4-oxo-7-(1-piperazinyl)quinoline-3-carboxylic acid (I) was synthesized from 2,4-dichloro-5-fluoroacetophenone through β-keto-ester formation, condensation with tri-Et orthoformate, substitution with 2-aminopyridine, cyclization, chelation with boric acid in acetic anhydride and followed by nucleophilic substitution reaction with piperazine. The total yield was 39.3%. The in vitro antibacterial activity of I against S. aureus and E. coli was tested.
IT **717910-04-8P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(synthesis and antibacterial activity of N-pyridine quinolone derivative)
RN 717910-04-8 CAPLUS
CN Boron, bis(acetato-κO)(7-chloro-6-fluoro-1,4-dihydro-4-(oxo-κO)-1-(2-pyridinyl)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)



L12 ANSWER 11 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2004:39482 CAPLUS
DOCUMENT NUMBER: 140:94138
TITLE: Ionic liquids containing borate or phosphate anions
INVENTOR(S): Moulton, Roger
PATENT ASSIGNEE(S): USA
SOURCE: U.S. Pat. Appl. Publ., 10 pp.
CODEN: USXXCO
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 2004007693	A1	20040115	US 2002-188452	20020703
WO 2004005222	A2	20040115	WO 2003-US21125	20030703
WO 2004005222	A3	20040923		

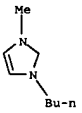
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

PRIORITY APPLN. INFO.: US 2002-188452 A 20020703

OTHER SOURCE(S): CASREACT 140:94138; MARPAT 140:94138
AB The present invention relates to novel ionic liqs. comprising a phosphate or borate anion. The ionic liqs. were prepared via metathesis or via a reaction between boric or phosphoric acid with metal hydroxide and an alc.
Thus, reaction of LiOH with boric acid in the presence of 3-methylsalicylic acid in water gave 70% lithium bis(3-methylsalicylyl)borate which on treatment with 1-butyl-3-methylimidazolium chloride in water gave 50% title ionic liquid, 1-butyl-3-methylimidazolium bis(3-methylsalicylyl)borate.
IT **625835-88-3P 642485-74-3P 642485-89-0P**
642485-99-2P 642486-04-2P 642486-27-9P
642486-33-7P 642486-38-2P
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of ionic liqs. containing borate or phosphate anions)
RN 625835-88-3 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[2-(hydroxy-κO)benzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

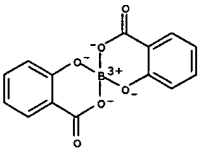
CRN 80432-08-2
CMF C8 H15 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

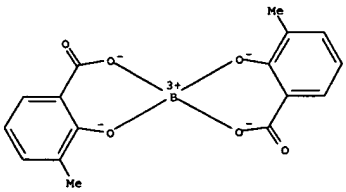
CRN 38403-08-6
CMF C14 H8 B O6
CCI CCS



RN 642485-74-3 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[2-(hydroxy-κO)-3-methylbenzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

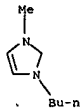
CRN 258875-08-0
CMF C16 H12 B O6
CCI CCS



L12 ANSWER 11 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CM 2

CRN 80432-08-2
CMF C8 H15 N2

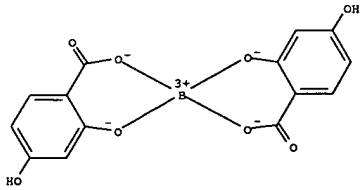


ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

RN 642485-89-0 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[2-(hydroxy-κO)-4-hydroxybenzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

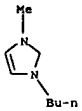
CM 1

CRN 642485-88-9
CMF C14 H8 B O8
CCI CCS



CM 2

CRN 80432-08-2
CMF C8 H15 N2

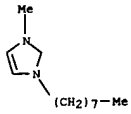


ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

RN 642485-99-2 CAPLUS
CN 1H-Imidazolium, 1-methyl-3-octyl-, (T-4)-bis[2-(hydroxy-κO)benzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

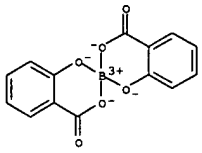
CRN 178631-03-3
CMF C12 H23 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

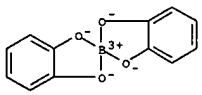
CRN 38403-08-6
CMF C14 H8 B O6
CCI CCS



RN 642486-04-2 CAPLUS
CN 1H-Imidazolium, 1-dodecyl-3-methyl-, (T-4)-bis[2-(hydroxy-κO)benzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

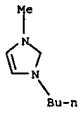
CRN 46928-10-3
CMF C16 H31 N2



RN 642486-33-7 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[1,2-benzenediolato(2-)-κO,κO']borate(1-) (9CI) (CA INDEX NAME)

CM 1

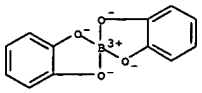
CRN 80432-08-2
CMF C8 H15 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

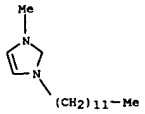
CRN 16986-25-7
CMF C12 H8 B O4
CCI CCS



RN 642486-38-2 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[4-(1,1-dimethylethyl)-1,2-benzenediolato(2-)-κO,κO']borate(1-) (9CI) (CA INDEX NAME)

CM 1

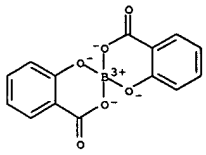
CRN 80432-08-2
CMF C8 H15 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

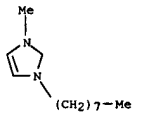
CRN 38403-08-6
CMF C14 H8 B O6
CCI CCS



RN 642486-27-9 CAPLUS
CN 1H-Imidazolium, 1-methyl-3-octyl-, (T-4)-bis[1,2-benzenediolato(2-)-κO,κO']borate(1-) (9CI) (CA INDEX NAME)

CM 1

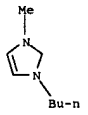
CRN 178631-03-3
CMF C12 H23 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

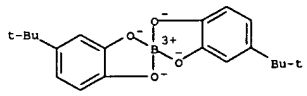
CRN 16986-25-7
CMF C12 H8 B O4
CCI CCS



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

CRN 53992-96-4
CMF C20 H24 B O4
CCI CCS



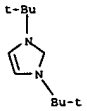
L12 ANSWER 12 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2003:892313 CAPLUS
DOCUMENT NUMBER: 139:352743
TITLE: Primary lithium batteries
INVENTOR(S): Munshi, M. Zafar A.; Coowar, Fazlil
PATENT ASSIGNEE(S): Lithium Power Technologies, Inc., USA
SOURCE: U.S. Pat. Appl. Publ., 6 pp.
CODEN: USXXCO
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 2003211383	A1	20031113	US 2002-142266	20020509
PRIORITY APPLN. INFO.:			US 2002-142266	20020509

AB A thermal battery for operation at temps. below about 250° and preferably not above about 200° includes a primarily CFx cathode, an electrolyte, and a lithium-based anode. The electrolyte is an organo borate lithium salt or an ionically conductive solid polymer electrolyte.
IT **618446-94-9**
RL: DEV (Device component use); USES (Uses)
(primary lithium thermal batteries)
RN 618446-94-9 CAPLUS
CN 1H-Imidazolium, 1,3-bis(1,1-dimethylethyl)-, (T-4)-bis[ethanedioato(2-)-κO1,κO2]borate(1-) (9CI) (CA INDEX NAME)

CM 1

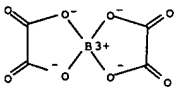
CRN 199382-55-3
CMF C11 H21 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

CRN 125579-65-9
CMF C4 B O8
CCI CCS

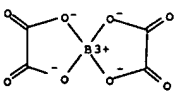


L12 ANSWER 12 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

L12 ANSWER 13 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2003:758951 CAPLUS
DOCUMENT NUMBER: 139:400500
TITLE: Ionic Liquids of Chelated Orthoborates as Model Ionic Glassformers
AUTHOR(S): Xu, Wu; Wang, Li-Min; Nieman, Ronald A.; Angell, C. Austen
CORPORATE SOURCE: Department of Chemistry and Biochemistry, Arizona State University, Tempe, AZ, 85287-1604, USA
SOURCE: Journal of Physical Chemistry B (2003), 107(42), 11749-11756
CODEN: JPCBFK; ISSN: 1520-6106
PUBLISHER: American Chemical Society
DOCUMENT TYPE: Journal
LANGUAGE: English
AB Ionic liqs. based on various chelated orthoborate anions of different N-containing onium cations have been synthesized using an economic synthesis strategy. Most orthoborates do not crystallize. They are found to have much higher glass transition temps. and room-temperature viscosities than those with perfluorinated anions such as TFSI-, BF-4, and CF3SO-3 (Tf-), as predicted from anion polarizability arguments. The ambient conductivities of the new ionic liqs. are low relative to those with perfluorinated anions. The transport properties all show that cohesion in these liqs. increases, and ionic mobilities decrease, as anion size increases, implying that van der Waals interactions, not Coulomb interactions, have become the controlling influence. In view of their resistance to crystallization, the large range of temperature over which these liqs. can be studied, their hydrophobic properties, and their high fragilities, these liqs. may provide good model systems for fundamental liquid state investigations and interesting solvents for large-mol. dissoln.
IT **566135-35-1P 625835-87-2P 625835-88-3P**
625835-89-4P 625835-90-7P 625835-91-8P
RL: PEP (Physical, engineering or chemical process); PRP (Properties); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)
(ionic liqs. of chelated orthoborates as model ionic glass formers)
RN 566135-35-1 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[ethanedioato(2-)-κO1,κO2]borate(1-) (9CI) (CA INDEX NAME)

CM 1

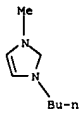
CRN 125579-65-9
CMF C4 B O8
CCI CCS



L12 ANSWER 13 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CM 2

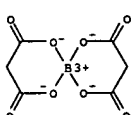
CRN 80432-08-2
CMF C8 H15 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE
RN 625835-87-2 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[propanedioato(2-)-κO1,κO3]borate(1-) (9CI) (CA INDEX NAME)

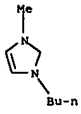
CM 1

CRN 343783-57-3
CMF C6 H4 B O8
CCI CCS



CM 2

CRN 80432-08-2
CMF C8 H15 N2

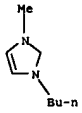


ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE
RN 625835-88-3 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[2-(hydroxy-

L12 ANSWER 13 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
κO}benzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

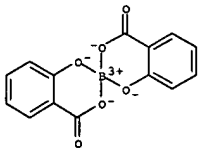
CRN 80432-08-2
CMF C8 H15 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

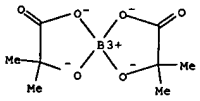
CRN 38403-08-6
CMF C14 H8 B O6
CCI CCS



RN 625835-89-4 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[2-(hydroxy-κO)-2-methylpropanoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

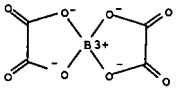
CRN 125579-63-7
CMF C8 H12 B O6
CCI CCS



L12 ANSWER 13 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CM 2

CRN 125579-65-9
CMF C4 B O8
CCI CCS

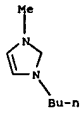


REFERENCE COUNT: 38 THERE ARE 38 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE
FORMAT

L12 ANSWER 13 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CM 2

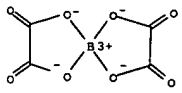
CRN 80432-08-2
CMF C8 H15 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE
RN 625835-90-7 CAPLUS
CN Pyridinium, 1-butyl-, (T-4)-bis[ethanedioato(2-)-κO1,κO2]borate(1-) (9CI) (CA INDEX NAME)

CM 1

CRN 125579-65-9
CMF C4 B O8
CCI CCS



CM 2

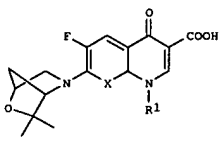
CRN 45806-95-9
CMF C9 H14 N

RN 625835-91-8 CAPLUS
CN Pyrrolidinium, 1-butyl-1-methyl-, (T-4)-bis[ethanedioato(2-)-κO1,κO2]borate(1-) (9CI) (CA INDEX NAME)

CM 1

CRN 223437-10-3
CMF C9 H20 N

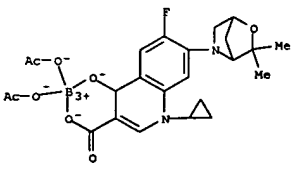
L12 ANSWER 14 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2003:557902 CAPLUS
DOCUMENT NUMBER: 140:217609
TITLE: Synthesis and antibacterial activity of 7-[(1S,4S)-3,3-dimethyl-2-oxa-5-azabicyclo[2.2.1]heptane-5-yl]quinolones
AUTHOR(S): Liu, Kaixiang; Guo, Huiyuan
CORPORATE SOURCE: Institute of Medicinal Biotechnology, Chinese Academy of Medical Sciences and Peking Union Medical College, Beijing, 100050, Peop. Rep. China
SOURCE: Zhongguo Yiyao Gongye Zazhi (2002), 33(6), 261-265
CODEN: ZYGZEA; ISSN: 1001-8255
PUBLISHER: Zhongguo Yiyao Gongye Zazhi Bianjibu
DOCUMENT TYPE: Journal
LANGUAGE: Chinese
OTHER SOURCE(S): CASREACT 140:217609
GI



I

II

AB Three Title compds. I (R1 = cyclopropyl, 2,4-difluorophenyl; X = C, N) were synthesized from trans-4-hydroxy-L-proline in eight steps to obtain an intermediate II, further substitution reaction of quinolone derivs., provide the title products. Their MICs against sixteen clin. isolates were detected. The results showed that they had very low antibacterial activities compared to ciprofloxacin and gatifloxacin.
IT **663605-48-9P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(synthesis of dimethyloxaazabicycloheptanquinolones)
RN 663605-48-9 CAPLUS
CN Borate(1-), bis(acetato-κO)[1-cyclopropyl-7-[(1S,4S)-3,3-dimethyl-2-oxa-5-azabicyclo[2.2.1]hept-5-yl]-6-fluoro-1,4-dihydro-4-(hydroxy-κO)-3-quinolinecarboxylato(2-)-κO3]-, (T-4)- (9CI) (CA INDEX NAME)



L12 ANSWER 14 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

L12 ANSWER 15 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:418223 CAPLUS
DOCUMENT NUMBER: 139:139046
TITLE: Ionic Liquids: Ion Mobilities, Glass Temperatures, and Fragilities
AUTHOR(S): Xu, Wu; Cooper, Emanuel I.; Angell, C. Austen
CORPORATE SOURCE: Department of Chemistry and Biochemistry, Arizona State University, Tempe, AZ, 85287-1604, USA
SOURCE: Journal of Physical Chemistry B (2003), 107(25), 6170-6178
CODEN: JPCBFK; ISSN: 1520-6106
PUBLISHER: American Chemical Society
DOCUMENT TYPE: Journal
LANGUAGE: English

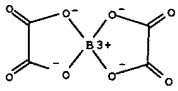
AB We combine old, unpublished data on ionic liqs. containing quaternary ammonium cations with new data on salts of aromatic cations containing a variety of anions, to demonstrate the existence for ionic liqs. of an unexpectedly wide range of liquid fragilities. The pattern is one now familiar for other liqs. Here, the pattern is important in determining the relative fluid properties at ambient temps. We find that the optimization of ionic liqs. for ambient temperature applications requiring low-vapor-pressure fluid phases involves the proper interplay of both cohesive energy and fragility factors. The cohesive energy is discussed in terms of the coulomb and van der Waals contributions to the attractive part of the pair potential. On the basis of the relation between the glass-transition temperature and the molar volume for salts with less-polarizable anions, we find evidence for a broad min. in the ionic liquid cohesive energy at an internuclear separation of ca. 0.6 nm. This min. lies between those of the BF4- and TFSI- anions for the small quaternary ammonium cations of this study. The min. is expected to be narrower and less well-defined for salts with polarizable anions. The relation of fluidity to conductance is considered in terms of a Walden plot that is shown to provide a useful basis for organizing the observations on ionic liqs. and solns. Low vapor pressure and ideal Walden products are intimately related.

IT **566135-35-1**
RL: PRP (Properties)
(ion mobilities, glass temps., and fragilities of ionic liqs.)

RN 566135-35-1 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, (T-4)-bis[ethanedioato(2-)-κO1,κO2]borate(1-) (9CI) (CA INDEX NAME)

CM 1

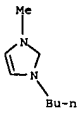
CRN 125579-65-9
CMF C4 B O8
CCI CCS



L12 ANSWER 15 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CM 2

CRN 80432-08-2
CMF C8 H15 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE
REFERENCE COUNT: 68 THERE ARE 68 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

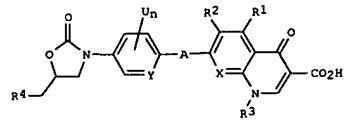
L12 ANSWER 16 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:319692 CAPLUS
DOCUMENT NUMBER: 138:338143
TITLE: Preparation of dual action bactericides comprising a oxazolidinone and a quinolone or naphthyridinone moiety effective against multi-drug resistant bacteria
INVENTOR(S): Hubschwerlen, Christian; Specklin, Jean-Luc
PATENT ASSIGNEE(S): Morphochem Aktiengesellschaft fuer Kombinatorische Chemie, Germany
SOURCE: PCT Int. Appl., 101 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 2
PATENT INFORMATION:

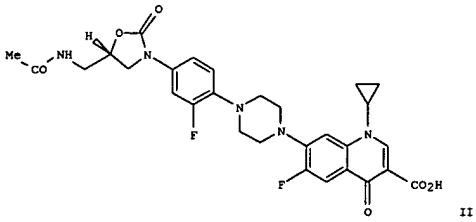
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003032962	A2	20030424	WO 2002-EP11163	20021004
WO 2003032962	A3	20030717		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW				
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
CA 2460572	AA	20030424	CA 2002-2460572	20021004
EP 1432705	A2	20040630	EP 2002-796533	20021004
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, SK				
BR 2002013063	A	20040928	BR 2002-13063	20021004
US 2005096343	A1	20050505	US 2003-491519	20021004
JP 2005529061	T2	20050929	JP 2003-535766	20021004
NZ 531879	A	20051028	NZ 2002-531879	20021004
ZA 2004001909	A	20050309	ZA 2004-1909	20040309
PRIORITY APPLN. INFO.:			US 2001-327162P	P 20011004
			WO 2002-EP11163	W 20021004

OTHER SOURCE(S): MARPAT 138:338143
GI

L12 ANSWER 16 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



I



II

AB The present invention relates to compds. of the Formula (I) that are useful antimicrobial agents and effective against a variety of multi-drug resistant bacteria. The present invention relates to oxazolidinones having a quinolone or naphthyridinone moiety (shown as I; variables defined below; e.g. 7-[4-[4-[(5S)-5-(acetylaminomethyl)-2-oxooxazolidin-3-yl]-2-fluorophenyl]piperazin-1-yl]-1-cyclopropyl-6-fluoro-4-oxo-1,4-dihydroquinoline-3-carboxylic acid (shown as II)) that are useful antibacterial agents and effective against a variety of multi-drug resistant bacteria. For I: A is a bond, NH, O, S, SO, SO2, SO2NH, PO4, -NH-CO-NH-, -CO-NH-, -CO-, -CO-O-, -NH-CO-O-, alkylene, alkenylene, alkynylene, heteroalkylene, arylene, heteroarylene, cycloalkylene, heterocycloalkylene, alkylarylene or heteroarylalkylene or a combination of two or more of these atoms or groups. X is CR5 or N; Y is CR6 or N; U is F or Cl; n = 0-3; R1 is H, F, Cl, Br, I, OH, NH2, alkyl or heteroalkyl; R2 is H, F or Cl; R3 is H, alkyl, alkenyl, alkynyl, heteroalkyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, alkylaryl or heteroarylalkyl; R4 is heteroalkyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, alkylaryl or heteroarylalkyl; R5 is H, F, Cl, OH, NH2, alkyl or heteroalkyl, or R3 and R5 can be linked via an alkylene, an alkenylene or heteroalkylene or be a part of a cycloalkylene or heterocycloalkylene group, in which case R3 is not H and R5 is not H, F, OH, NH2 or Cl; R6 is H, F, Cl or OMe. Although the methods of preparation are not claimed, 30 example prepns. are included; the examples of this patent and many of the claims are the same as those of WO 03/031443 A1. All examples were tested against several gram pos. and gram neg. bacteria; typical MIC ranges (mg/L) are: S. aureus (MRSA: 0.125-2; MSSA: 0.06-1), E. faecalis (≤0.03-1), E. faecium (≤0.03-1), and S. pneumoniae (≤0.03-1). They all have a broader and more pronounced activity

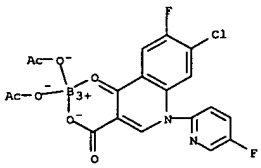
L12 ANSWER 16 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

than the corresponding quinolone and oxazolidinone as well as a 1+1 combination of these two compds.

IT **510728-72-0P**, 7-Chloro-6-fluoro-1-(5-fluoropyridin-2-yl)-4-oxo-1,4-dihydroquinoline-3-carboxylatoboron diacetate
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation of dual action bactericides comprising oxazolidinone and quinolone or naphthyridinone moiety effective against multi-drug resistant bacteria)

RN 510728-72-0 CAPLUS

CN Boron, bis(acetato-κO)[7-chloro-6-fluoro-1-(5-fluoro-2-pyridinyl)-1,4-dihydro-4-(oxo-κO)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)



L12 ANSWER 17 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:301084 CAPLUS
DOCUMENT NUMBER: 138:304289
TITLE: Preparation of dual action bactericides comprising a oxazolidinone and a quinolone or naphthyridinone moiety effective against multi-drug resistant bacteria
INVENTOR(S): Hubschwerlen, Christian; Specklin, Jean-Luc
PATENT ASSIGNEE(S): Morphochem Aktiengesellschaft fuer Kombinatorische Chemie, Germany
SOURCE: PCT Int. Appl., 100 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 2
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003031443	A1	20030417	WO 2002-EP10766	20020925
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW				
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
ZA 2004001909	A	20050309	ZA 2004-1909	20040309
PRIORITY APPLN. INFO.:			US 2001-327162P P	20011004

OTHER SOURCE(S): MARPAT 138:304289
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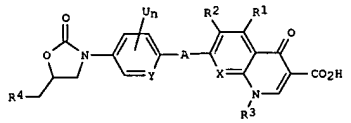
L12 ANSWER 17 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



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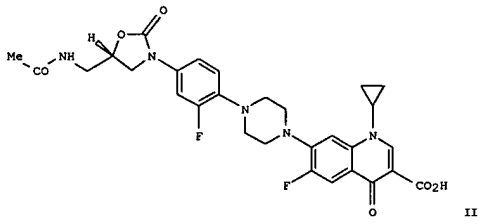
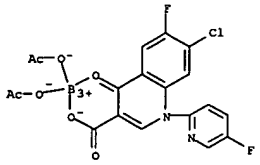


II

AB The present invention relates to oxazolidinones having a quinolone or naphthyridinone moiety (shown as I; variables defined below; e.g. 7-[4-[4-[(5S)-5-(acetylaminomethyl)-2-oxooxazolidin-3-yl]-2-fluorophenyl]piperazin-1-yl]-1-cyclopropyl-6-fluoro-4-oxo-1,4-dihydroquinoline-3-carboxylic acid (shown as II)) that are useful antibacterial agents and effective against a variety of multi-drug resistant bacteria. For I: A is a bond, NH, O, S, SO, SO2, SO2NH, PO4, -NH-CO-NH-, -CO-NH-, -CO-, -CO-O-, -NH-CO-O-, alkylene, alkenylene, alkynylene, heteroalkylene, arylene, heteroarylene, cycloalkylene, heterocycloalkylene, alkylarylene or heteroarylalkylene or a combination of two or more of these atoms or groups. X is CR5 or N; Y is CR6 or N; U is F or Cl; n = 0-3; R1 is H, F, Cl, Br, I, OH, NH2, alkyl or heteroalkyl; R2 is H, F or Cl; R3 is H, alkyl, alkenyl, alkynyl, heteroalkyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, alkylaryl or heteroarylalkyl; R4 is heteroalkyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, alkylaryl or heteroarylalkyl; R5 is H, F, Cl, OH, NH2, alkyl or heteroalkyl, or R3 and R5 can be linked via an alkylene, an alkenylene or heteroalkylene or be a part of a cycloalkylene or heterocycloalkylene group, in which case R3 is not H and R5 is not H, F, OH, NH2 or Cl; R6 is H, F, Cl or OMe. Although the methods of preparation are not claimed, 30 example prepns. are included. All examples were tested against several gram pos. and gram neg. bacteria; typical MIC ranges (mg/L) are: S. aureus (MRSA: 0.125-2; MSSA: 0.06-1), E. faecalis (≤0.03-1), E. faecium (≤0.03-1), and S. pneumoniae (≤0.03-1). They all have a broader and more pronounced activity than the corresponding quinolone and oxazolidinone as well as a 1+1 combination of these two compds.

IT **510728-72-0P**, 7-Chloro-6-fluoro-1-(5-fluoropyridin-2-yl)-4-oxo-1,4-dihydroquinoline-3-carboxylatoboron diacetate
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT

L12 ANSWER 17 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
(Reactant or reagent)
(prepn. of dual action bactericides comprising oxazolidinone and quinolone or naphthyridinone moiety effective against multi-drug resistant bacteria)
RN 510728-72-0 CAPLUS
CN Boron, bis(acetato-κO)[7-chloro-6-fluoro-1-(5-fluoro-2-pyridinyl)-1,4-dihydro-4-(oxo-κO)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE

FORMAT

L12 ANSWER 18 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2003:301082 CAPLUS
DOCUMENT NUMBER: 138:304288
TITLE: Preparation of dual action bactericides comprising a oxazolidinone and a quinolone or naphthyridinone moiety effective against multi-drug resistant bacteria
INVENTOR(S): Hubschwerlen, Christian; Specklin, Jean-Luc
PATENT ASSIGNEE(S): Morphochen Aktiengesellschaft fuer Kombinatorische Chemie, Germany
SOURCE: PCT Int. Appl., 95 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

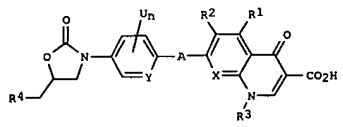
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003031441	A1	20030417	WO 2002-EP10765	20020925

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

PRIORITY APPLN. INFO.: US 2001-327208P P 20011004

OTHER SOURCE(S): MARPAT 138:304288
GI

L12 ANSWER 18 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



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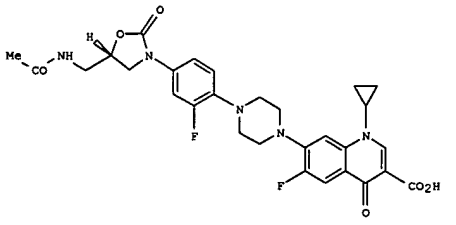
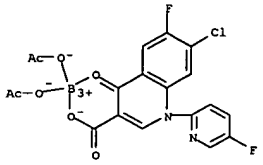


II

AB The present invention refers to novel multiple action compds., i.e., to compds. which contain at least two pharmaceutically active components in one mol. The compds. have a higher stability than corresponding compds. of the prior art. Although the present invention does not claim any specific compds. or even a Markush expression, the examples involve oxazolidinones having a quinolone or naphthyridinone moiety (shown as I; variables defined below; e.g. 7-[4-[4-[(5S)-5-(acetylaminomethyl)-2-oxooxazolidin-3-yl]-2-fluorophenyl]piperazin-1-yl]-1-cyclopropyl-6-fluoro-4-oxo-1,4-dihydroquinoline-3-carboxylic acid (shown as II)) that are useful antibacterial agents and effective against a variety of multi-drug resistant bacteria. For I: A is a bond, NH, O, S, SO, SO2, SO2NH, PO4, -NH-CO-NH-, -CO-NH-, -CO-, -CO-O-, -NH-CO-O-, alkylene, alkenylene, alkynylene, heteroalkylene, arylene, heteroarylene, cycloalkylene, heterocycloalkylene, alkylarylene or heteroarylalkylene or a combination of two or more of these atoms or groups. X is CR5 or N; Y is CR6 or N; U is F or Cl; n = 0-3; R1 is H, F, Cl, Br, I, OH, NH2, alkyl or heteroalkyl; R2 is H, F or Cl; R3 is H, alkyl, alkenyl, alkynyl, heteroalkyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, alkylaryl or heteroarylalkyl; R4 is heteroalkyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, alkylaryl or heteroarylalkyl; R5 is H, F, Cl, OH, NH2, alkyl or heteroalkyl, or R3 and R5 can be linked via an alkylene, an alkenylene or heteroalkylene or be a part of a cycloalkylene or heterocycloalkylene group, in which case R3 is not H and R5 is not H, F, OH, NH2 or Cl; R6 is H, F, Cl or OMe. Although the methods of preparation are not claimed, 30 example prepns. are included. All examples were tested against several gram pos. and gram neg. bacteria; typical MIC ranges (mg/L) are: S. aureus (MRSA: 0.125-2; MSSA: 0.06-1), E. faecalis (≤0.03-1), E. faecium (≤0.03-1), and S. pneumoniae (≤0.03-1). They all have a broader and more pronounced activity than the corresponding quinolone and oxazolidinone as well as a 1+1 combination of these two compds. The examples of this patent are the same as those of WO 03/031443 A1.

L12 ANSWER 18 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
IT 510728-72-0P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation of dual action bactericides comprising oxazolidinone and quinolone or naphthyridinone moiety effective against multi-drug resistant bacteria)
RN 510728-72-0 CAPLUS
CN Boron, bis(acetato-κO)[7-chloro-6-fluoro-1-(5-fluoro-2-pyridinyl)-1,4-dihydro-4-(oxo-κO)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE

FORMAT

L12 ANSWER 19 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:200758 CAPLUS
DOCUMENT NUMBER: 138:245327
TITLE: Organic electroluminescent device with boron tetraquinolinate derivative
INVENTOR(S): Suzuki, Koichi; Senoo, Akihiro; Sven, Anderssen; Ueno, Kazunori
PATENT ASSIGNEE(S): Canon Inc., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 23 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2003077671	A2	20030314	JP 2001-265872	20010903
PRIORITY APPLN. INFO.:			JP 2001-265872	20010903

OTHER SOURCE(S): MARPAT 138:245327

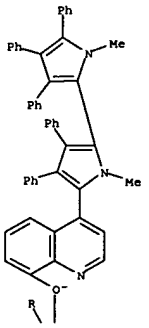
AB The invention refers to an organic electroluminescent device comprising a boron tetraquinolinate derivative (Markush structures provided) in the organic layer.

IT **501667-85-2 501667-86-3 501667-87-4 501667-88-5 501667-90-9 501667-92-1 501667-93-2 501667-98-7 501667-99-8 501668-09-3**
RL: DEV (Device component use); USES (Uses)
(organic electroluminescent device with boron tetraquinolinate derivative)

RN 501667-85-2 CAPLUS
CN Borate(1-), tetrakis[4-(1,1'-dimethyl-3,3',4,4',5'-pentaphenyl[2,2'-bi-1H-pyrrol]-5-yl)-8-quinolinolato-κO8]-, lithium (9CI) (CA INDEX NAME)

L12 ANSWER 19 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

PAGE 1-A



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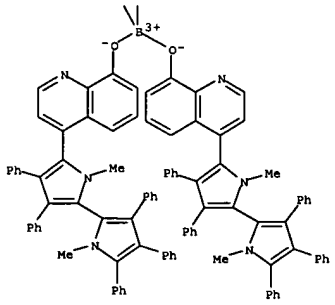
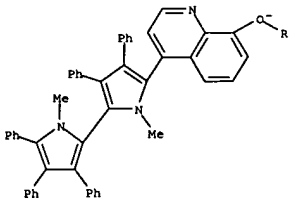


L12 ANSWER 19 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

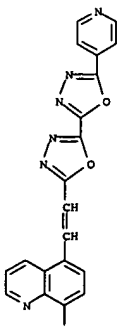
PAGE 3-A



● Li+

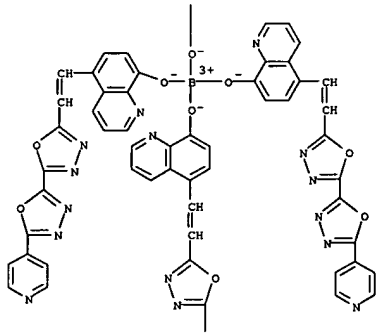
RN 501667-86-3 CAPLUS
CN Borate(1-), tetrakis[5-[2-[5'-(4-pyridinyl)[2,2'-bi-1,3,4-oxadiazol]-5-yl]ethenyl]-8-quinolinolato-κO8]-, potassium (9CI) (CA INDEX NAME)

PAGE 1-A



L12 ANSWER 19 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

PAGE 2-A



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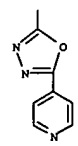
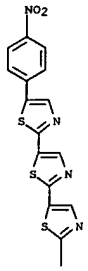
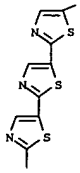


● K+

RN 501667-87-4 CAPLUS
CN Borate(1-), tetrakis[2-[2-[5-(4-nitrophenyl)[2,5':2',5'':2'',5''':2''',5''''-quinquethiazol]-2''''-yl]ethenyl]-8-quinolinolato-κO8]-, sodium (9CI) (CA INDEX NAME)

PAGE 1-A





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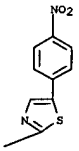

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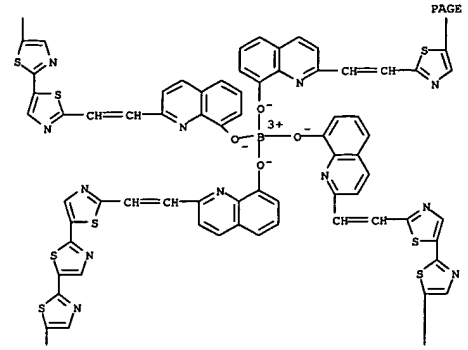

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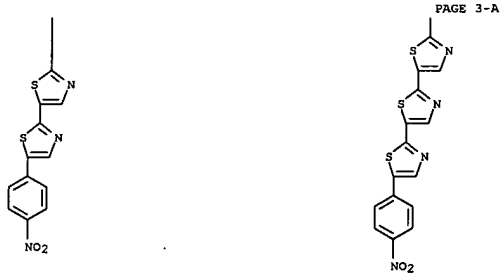
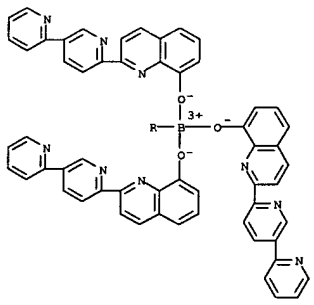


● Na+

RN 501667-88-5 CAPLUS
CN Borate(1-), tetrakis(2-[2,3'-bipyridin]-5'-yl-8-quinolinolato-κO8)-, lithium (9CI) (CA INDEX NAME)

PAGE 1-A



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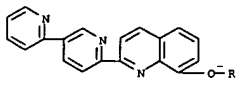
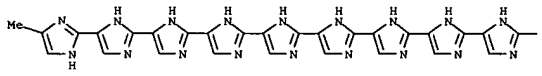


● Li+

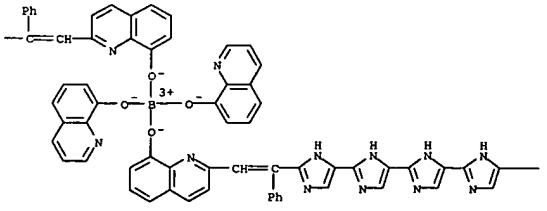
RN 501667-90-9 CAPLUS
CN Borate(1-), bis[2-[2-(4-methyl[2,4':2',4'':2'',4''':2''',4'''':2'''',4''''':2''''',4'''''':2'''''',4''''''':2''''''',4'''''''-novi-1H-imidazol]-2'''''''-yl)-2-phenylethenyl]-8-quinolinolato-κO8]bis(8-quinolinolato-κO8)-, lithium, (T-4)- (9CI) (CA INDEX NAME)

PAGE 1-A



● Li+

PAGE 1-B



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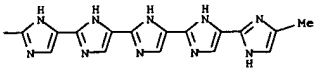
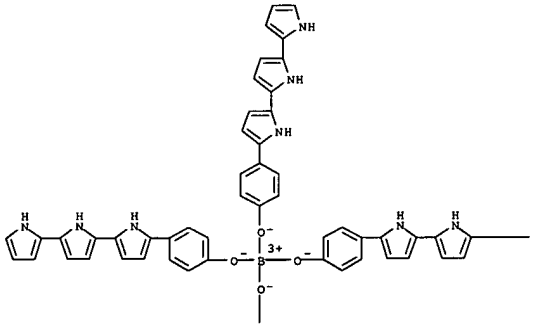


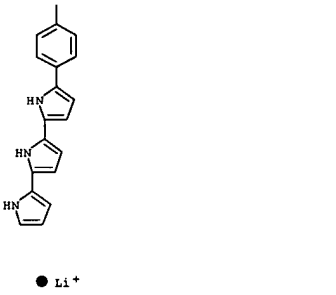
RN 501667-92-1 CAPLUS
CN Borate(1-), tetrakis(4-[2,2':5',2''-ter-1H-pyrrol]-5-ylphenolato-κO)-, lithium (9CI) (CA INDEX NAME)

PAGE 1-A



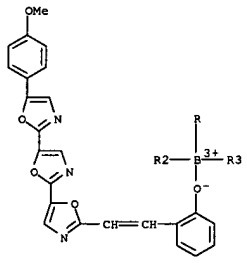
PAGE 1-B

PAGE 2-A



RN 501667-93-2 CAPLUS
CN Borate(1-), tetrakis[2-[2-[5-(4-methoxyphenyl)[2,5':2',5''-teroxazol]-2''-yl]ethenyl]phenolato-κO]-, potassium (9CI) (CA INDEX NAME)

PAGE 1-A



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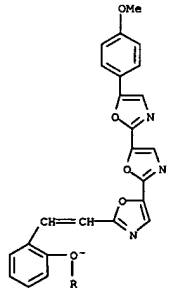

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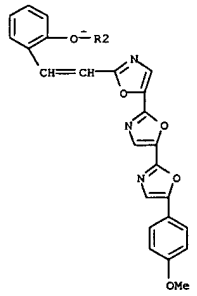

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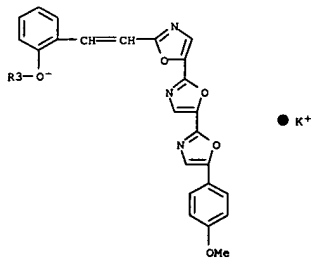

RN 501667-98-7 CAPLUS
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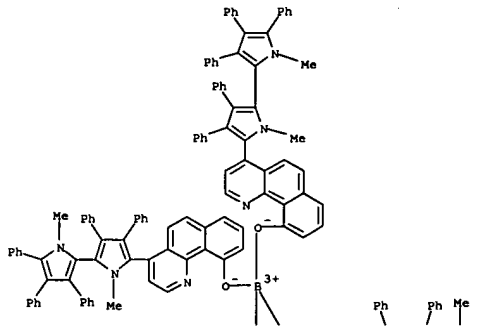
PAGE 1-A



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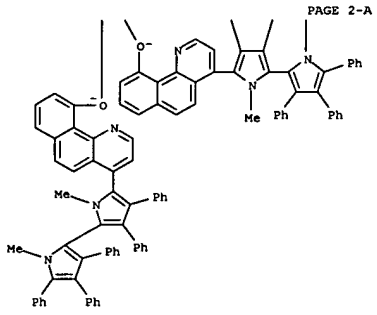
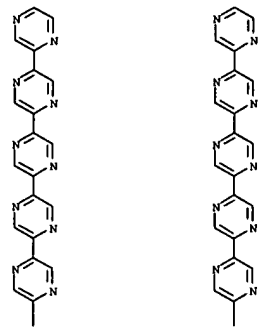


PAGE 3-A

RN 501667-99-8 CAPLUS
CN Borate(1-), tetrakis(2-[2,2':5',2'':5'',2''':5''',2'''':4'''',2''''':5'''''-sexipyrazin]-5-ylbenzo[h]quinolin-10-olato-κO10)-, lithium (9CI) (CA INDEX NAME)

PAGE 1-A



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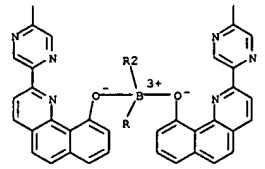

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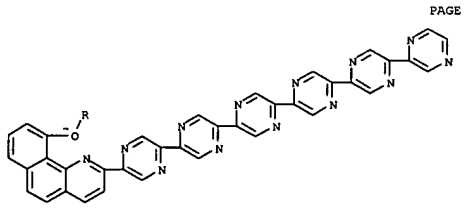

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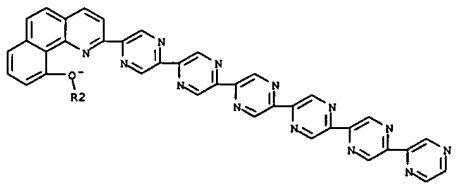

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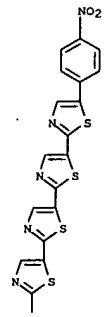
PAGE 1-A



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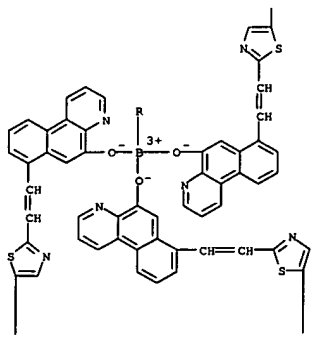

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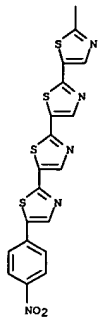

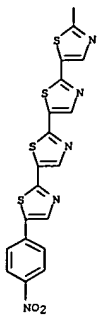

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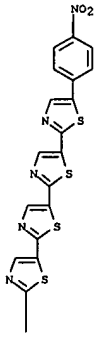

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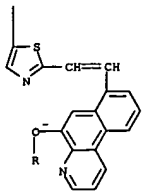


ACCESSION NUMBER: 2003:107142 CAPLUS
DOCUMENT NUMBER: 138:296637
TITLE: [Mo5VMo7VIO30(BPO4)2(O3P-Ph)6]5-: A Phenyl-Substituted Molybdenum(V/VI) Boro-Phosphate Polyoxometalate
AUTHOR(S): Sassoye, Capucine; Norton, Kieran; Sevov, Slavi C.
CORPORATE SOURCE: Department of Chemistry and Biochemistry, University of Notre Dame, Notre Dame, IN, 46556, USA
SOURCE: Inorganic Chemistry (2003), 42(5), 1652-1655
CODEN: INOCAJ; ISSN: 0020-1669
PUBLISHER: American Chemical Society
DOCUMENT TYPE: Journal
LANGUAGE: English
OTHER SOURCE(S): CASREACT 138:296637

AB The polyanion [Mo5VMo7VIO30(BPO4)2(O3PPh)6]5- is the first hybrid borophosphate-phenylphosphonate polyoxometalate. It was structurally characterized as its imidazolium salt, (C3N2H5)5[Mo12O30(BPO4)2(O3PPh)6].cntdot.H2O (monoclinic, P21/c, a 22.120(3), b 13.042(2), and c 32.632(4) Å, β 101.293(3)), which was synthesized hydrothermally from imidazole, molybdenum oxide and metal, and boric, phosphoric, and phenylphosphonic acids. The anion is the second example of a new class of polyoxometalates that resemble Dawson anions but where the two pole caps of three edge-sharing MoO6 octahedra in the latter are replaced by other units, in this case tetrahedral borate sharing corners with three phenylphosphonic groups, {(OB)(O3P-Ph)3}. The 12 molybdenum atoms forming the two equatorial belts of the cluster are of mixed-valence, five are MoV and seven are MoVI, and the resulting five electrons are delocalized. Four of these electrons are paired according to the temperature dependence of the magnetic susceptibility. The new compound is soluble in a mixture of water and pyridine (in equal vols.) as well as in nitromethane, and the anions are intact in these solns.

IT **506413-04-3P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of molybdenum borophosphate phenylphosphonate polyoxometalate)

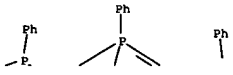
RN 506413-04-3 CAPLUS

CN Molybdate(5-), diborateoctadeca-μ-oxododecaoxohexakis[μ3-[phenylphosphonato(2-)-κO:κO':κO'']]bis[μ7-[phosphato(3-)-κO:κO:κO':κO':κO'':κO'':κO''']]dodeca-, pentahydrogen, compd. with pyridine (1:5) (9CI) (CA INDEX NAME)

CM 1

CRN 506413-02-1
CMF C36 H30 B2 Mo12 O56 P8 . 5 H
CCI CCS

PAGE 1-A



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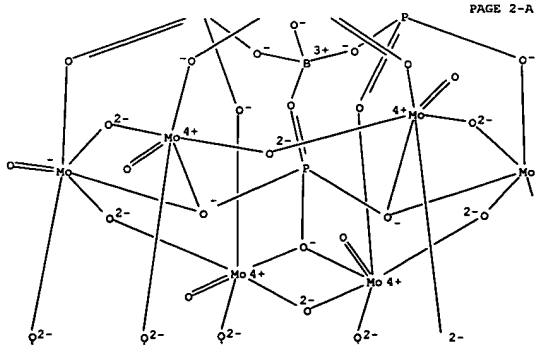

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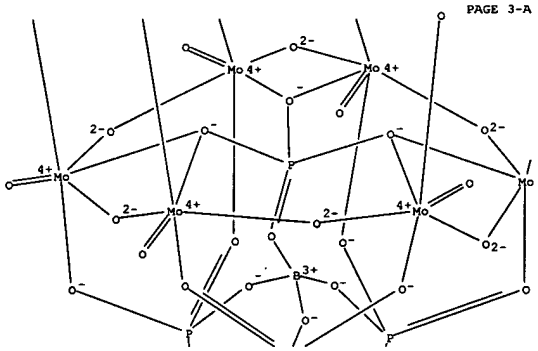

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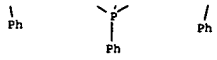
PAGE 4-A



●5 H^+

CM 2

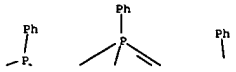
CRN 110-86-1
CMF C5 H5 N

IT **506413-03-2P**
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(preparation, crystal structure and magnetic susceptibility of molybdenum
borophosphate phenylphosphonate polyoxometalate)
RN 506413-03-2 CAPLUS
CN Molybdate(5-), octadeca-μ-oxododecaoxohexakis[μ3-[phenylphosphonato(2-)-κO:κO':κO'']]bis[μ7-[phosphato(3-)-κO:κO:κO':κO':κO'':κO'' :κO''']]diboratedodeca-, pentahydrogen, compd. with 1H-imidazole (1:5), monohydrate (9CI) (CA INDEX NAME)

CM 1

CRN 506413-02-1
CMF C36 H30 B2 Mo12 O56 P8 . 5 H
CCI CCS

PAGE 1-A



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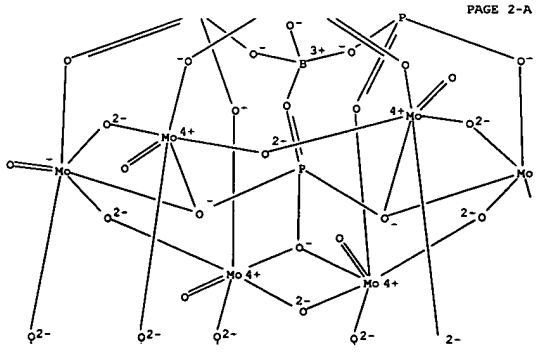

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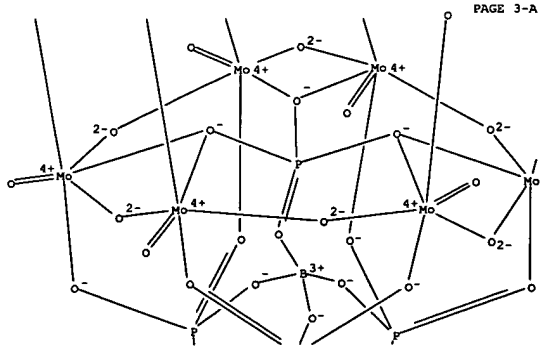
PAGE 3-A



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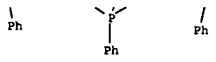
PAGE 4-A



●5 H+

CM 2

CRN 288-32-4
CMF C3 H4 N2

REFERENCE COUNT: 17 THERE ARE 17 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE
FORMAT

L12 ANSWER 21 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:22874 CAPLUS
DOCUMENT NUMBER: 138:89799
TITLE: Preparation of fluoroquinolonyl derivatives of oxazolidinones as antibacterial agents
INVENTOR(S): Mourelle Mancini, Marisabel; Huguet Clotet, Juan; Hidalgo Rodriguez, Jose; Del Castillo, Juan Carlos
PATENT ASSIGNEE(S): Vita-Invest, S.A., Spain
SOURCE: PCT Int. Appl., 110 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003002560	A1	20030109	WO 2002-IB2408	20020624
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO				
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
ES 2186550	A1	20030501	ES 2001-1559	20010627
ES 2186550	B2	20031116		
CA 2450982	AA	20030109	CA 2002-2450982	20020624
EE 200400004	A	20040216	EE 2004-4	20020624
EP 1401834	A1	20040331	EP 2002-738497	20020624
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JP 2004521147	T2	20040715	JP 2003-508941	20020624
CN 1520412	A	20040811	CN 2002-812852	20020624
BR 2002010667	A	20041005	BR 2002-10667	20020624
NZ 530206	A	20050729	NZ 2002-530206	20020624
US 2004147545	A1	20040729	US 2003-469283	20030828
NO 2003005791	A	20040219	NO 2003-5791	20031222
BG 108498	A	20050331	BG 2003-108498	20031222
PRIORITY APPLN. INFO.:			ES 2001-1559 A	20010627
			WO 2002-IB2408 W	20020624

OTHER SOURCE(S): MARPAT 138:89799
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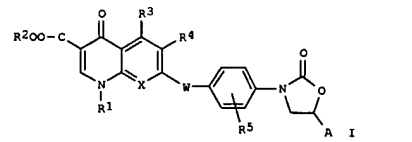

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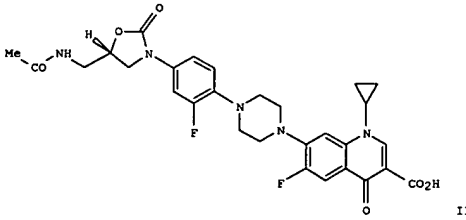


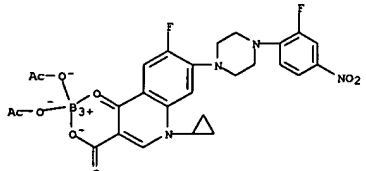
II

AB This invention discloses new fluoroquinolonyl derivs. of oxazolidinones (shown as I; variables defined below; e.g. 7-[4-[4-[5-(S)-(acetylaminomethyl)-2-oxooxazolidin-3-yl]-2-fluorophenyl]piperazin-1-yl]-1-cyclopropyl-6-fluoro-4-oxo-1,4-dihydroquinoline-3-carboxylic acid (shown as II)) and processes for obtaining them, the corresponding pharmaceutical compns. and use thereof for manufacturing a medicament for the treatment of microbial infections. These new compds. are useful as antibacterial agents. Furthermore phenalen-type compds. according to (II) are disclosed. Compds. I show activity as antibacterial agents; MIC values for .apprx.15 compds. are included. Advantageously they possess a broad spectrum of activity against gram-pos. bacteria such as Staphylococcus, Streptococcus, Enterococcus and the like, as well as against gram-neg. bacteria such as E. Coli, H. Influenzae, M. Catarrahalis, etc., and even against strains resistant to known antibiotics such as meticillin, vancomicin, penicillin, etc. They are also active against anaerobic microorganisms such as Bacteroides fragilis. Thirty-five example prepns. of I plus 38 example prepns. of intermediates are included. II was prepared from 7-[4-[-[5-(S)-(acetylaminomethyl)-2-oxooxazolidin-3-yl]-2-fluorophenyl]piperazin-1-yl]-1-cyclopropyl-6-fluoro-4-oxo-1,4-dihydroquinoline-3-carboxylic acid diacetoxyboron chelate in H2O and MeCN using 1N NaOH at room temperature; the chelate was prepared from N-[[3-(3-fluoro-4-(piperazin-1-yl)phenyl]-2-oxooxazolidin-5-(S)-yl]methyl]acetamide, 7-chloro-1-cyclopropyl-6-fluoro-4-oxo-1,4-dihydroquinoline-3-carboxylic acid diacetoxyboron chelate and Et3N in MeCN at reflux for 16 h. For I: X = CR6 or N; R1 = C1-C4-alkyl,

C3-C6-cycloalkyl, C2-C4-alkenyl, 2-hydroxyethyl, 2-fluoroethyl, or Ph optionally substituted by 1 or 2 atoms of F; R2: H, alkyl C1-C4 or phenyl; R3 = H, halogen, C1-C4-alkyl or C1-C4-alkoxy, amino; R4 = H or halogen; R6 = H, halogen, C1-C4-alkyl, C1-C4-haloalkoxy or else R1 and R6 together form a bridge of structure -CHMe-CH2-O-, -CHMe-CH2-S-, -CHMe-CH2-CH2-. R5 = H, halogen, OCH3, C1-C4-alkoxy, C1-C4-alkyl or C1-C4-haloalkyl; A = -CH2-NH-R7, -CHOH-C.tplbond.CH; wherein R7 = isoxazol, -CO-R8, -CS-R8, -CS-OR8, -COOR8, -CONHR8, -CSNHR8, -SO2R8 or COCH:CHAr (Ar = R9-substituted phenyl) wherein R8 = C1-C4-alkyl, C1-C4-haloalkyl, C2-C4-alkenyl, aryl, C1-C4-alkyl substituted by an C1-C4-alkoxy group, C1-C4-carboxyalkyl, cyano, or amino. R9= H, C1-C4-alkyl, C2-C4-alkenyl, OH, C1-C4-alkoxy, NR12R13, NO2, halogen, or CO-R12; R12 and R13 = H or C1-C4-alkyl; W = azetidinyl, pyrrolidinyl, azepanyl, and piperazinyl derivs. as more fully described in the claims.

IT **484638-76-8P 484638-77-9P 484639-06-7P 484639-07-8P 484639-08-9P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation of fluoroquinolonyl derivs. of oxazolidinones as antibacterial agents)

RN 484638-76-8 CAPLUS
CN Boron, bis(acetato-κO)[1-cyclopropyl-6-fluoro-7-[4-(2-fluoro-4-nitrophenyl)-1-piperazinyl]-1,4-dihydro-4-(oxo-κO)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)



RN 484638-77-9 CAPLUS
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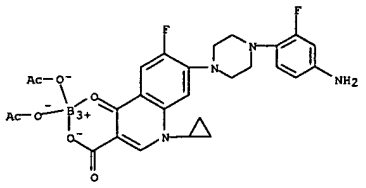
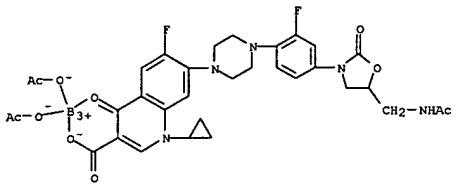
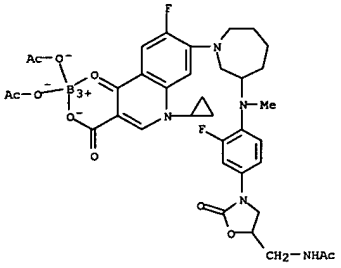


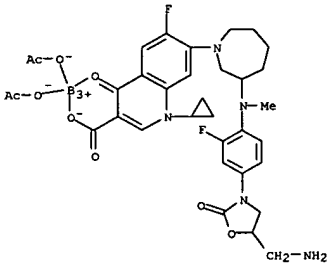
RN 484639-06-7 CAPLUS
CN Boron, bis(acetato-κO)[7-[4-[4-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-fluorophenyl]-1-piperazinyl]-1-cyclopropyl-6-fluoro-1,4-dihydro-4-(oxo-κO)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)



RN 484639-07-8 CAPLUS
CN Boron, bis(acetato-κO)[7-[3-[[4-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-fluorophenyl]methylamino]hexahydro-1H-azepin-1-yl]-1-cyclopropyl-6-fluoro-1,4-dihydro-4-(oxo-κO)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)



RN 484639-08-9 CAPLUS
CN Boron, bis(acetato-κO)[7-[3-[[4-[5-(aminomethyl)-2-oxo-3-oxazolidinyl]-2-fluorophenyl]methylamino]hexahydro-1H-azepin-1-yl]-1-cyclopropyl-6-fluoro-1,4-dihydro-4-(oxo-κO)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)

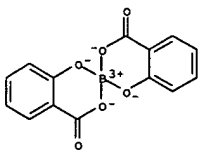


REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 22 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2002:789169 CAPLUS
DOCUMENT NUMBER: 139:16513
TITLE: Synthesis of bis(salicylato)borates, bis(naphthalene-2,3-diolato)borates and their analog
AUTHOR(S): Turiashvili, L. G.; Machkhoshvili, R. I.; Dzhioshvili, B. D.; Doksopulo, E. P.; Parunashvili, N. A.
CORPORATE SOURCE: Tbilis. Gos. Univ. im. Iv. Dzhavakhishvili, Tbilissi, Georgia
SOURCE: Sakartvelos Mecnierebata Akademiis Macne, Kimiis Seria (2002), 28(1-2), 45-51
CODEN: IANKEJ
PUBLISHER: Proizvodstvenno-Izdatel'skoe Ob'edinenie "Metsniereba"
DOCUMENT TYPE: Journal
LANGUAGE: Russian
OTHER SOURCE(S): CASREACT 139:16513
AB HL[BL1L2].nH2O (L = ethanolamine, ethylenediamine, N-phenylethylenediamine, pyridine, 2-amino-5-methylpyridine; H2L1 = H2L2 = salicylic acid, naphthalene-2,3-diol or HL1 = salicylic acid and H2L2 = naphthalene-2,3-diol) were prepared from B(OH)3, H2L1 and L. The composition and structure of the synthesized compds. were established by IR spectroscopy, thermal and chemical anal. and from elec. conductivity
IT **69030-98-4P 532927-80-3P 532927-81-4P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation and thermal decomposition)
RN 69030-98-4 CAPLUS
CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS



● H+

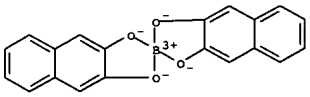
CM 2

CRN 110-86-1
CMF C5 H5 N

RN 532927-80-3 CAPLUS
CN Borate(1-), bis[2,3-naphthalenediolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with pyridine (1:1), dihydrate (9CI) (CA INDEX NAME)

CM 1

CRN 47422-29-7
CMF C20 H12 B O4 . H
CCI CCS



● H+

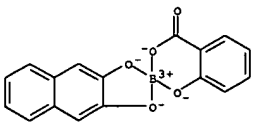
CM 2

CRN 110-86-1
CMF C5 H5 N

RN 532927-81-4 CAPLUS
CN Borate(1-), [2-(hydroxy-κO)benzoato(2-)-κO][2,3-naphthalenediolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with pyridine (1:1), monohydrate (9CI) (CA INDEX NAME)

CM 1

CRN 532927-77-8
CMF C17 H10 B O5 . H
CCI CCS



● H+

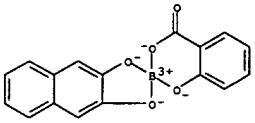
CM 2

CRN 110-86-1
CMF C5 H5 N

IT **532927-83-6P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 532927-83-6 CAPLUS
CN Borate(1-), [2-(hydroxy-κO)benzoato(2-)-κO][2,3-naphthalenediolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with 5-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

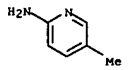
CRN 532927-77-8
CMF C17 H10 B O5 . H
CCI CCS



● H+

CM 2

CRN 1603-41-4
CMF C6 H8 N2

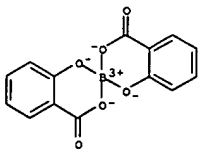


ACCESSION NUMBER: 2002:740867 CAPLUS
DOCUMENT NUMBER: 138:31671
TITLE: Development of high-temperature resistant electrolytic solutions
AUTHOR(S): Kikuchi, Kazunobu
CORPORATE SOURCE: 1st Dev. Technol. Div., Tomiyama Pure Chem. Ind., Ltd., Japan
SOURCE: Denkai Chikudenki Hyoron (2002), 53(1), 101-110
CODEN: DCHYAK; ISSN: 0286-5629
PUBLISHER: Denkai Chikudenki Kenkyukai
DOCUMENT TYPE: Journal
LANGUAGE: Japanese
AB Various materials were examined for high-temperature resisting electrolytic solns. by keeping at 120° for 100-1000 h to determine the elec. conductivity, pH, and chromatog. for various cation mixts. Boric acid-disalicylate was more temperature resisting than phthalates which produce impurities during heat treatment. The superior cations were amidine and cyclic tertiary amine.

IT **221332-52-1 478012-80-5 478012-81-6 478014-56-1**
RL: PRP (Properties); TEM (Technical or engineered material use); USES (Uses)
(development of high-temperature resistant electrolytic solns.)
RN 221332-52-1 CAPLUS
CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, (T-4)-, hydrogen, compd. with 1,2-dimethyl-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS



● H+

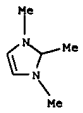
CM 2

CRN 1739-84-0
CMF C5 H8 N2

RN 478012-80-5 CAPLUS
CN 1H-Imidazolium, 1,2,3-trimethyl-, (T-4)-bis[2-(hydroxy-κO)benzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

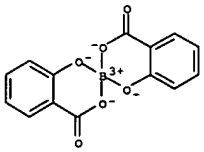
CRN 65086-10-4
CMF C6 H11 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

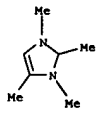
CRN 38403-08-6
CMF C14 H8 B O6
CCI CCS



RN 478012-81-6 CAPLUS
CN 1H-Imidazolium, 1,2,3,4-tetramethyl-, (T-4)-bis[2-(hydroxy-κO)benzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

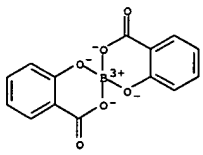
CRN 186612-73-7
CMF C7 H13 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

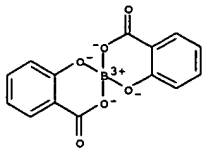
CRN 38403-08-6
CMF C14 H8 B O6
CCI CCS



RN 478014-56-1 CAPLUS
CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, (T-4)-, hydrogen, compd. with 1,4,5,6-tetrahydro-1,2-dimethylpyrimidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS



CM 2

CRN 4271-96-9
CMF C6 H12 N2

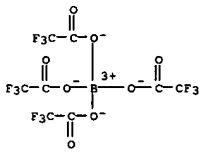
ACCESSION NUMBER: 2002:441202 CAPLUS
DOCUMENT NUMBER: 137:27396
TITLE: Preparation of tetrakis(acyloxy)borate(1-) and their substituted onium salts
INVENTOR(S): Nagata, Hiroshi
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 10 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2002167390	A2	20020611	JP 2000-364192	20001130
PRIORITY APPLN. INFO.:			JP 2000-364192	20001130

OTHER SOURCE(S): CASREACT 137:27396; MARPAT 137:27396
AB The title [A][(R1CO2)4B] (I; A = R2R3R4HN+, R5R6C:N+HR7; R1 = alkyl, aryl, aralkyl; R2-7 = alkyl, aryl; ≤2 of R2-R4 or R5-R7 may be bonded together to form a ring) are prepared by treating B(OH)3 with (R1CO)2O (R2 = same as above) and NR2R3R4 (R2-R4 = same as above) or R5R6C:NR7 (R5-R7 = same as above). Then, substituted onium salts of I, useful as curing accelerators for epoxy resins used in elec. and electronic materials, are prepared by treating I with tetrasubstituted phosphonium, tetrasubstituted ammonium, or trisubstituted sulfonium salts. E.g., H3BO3, 1-naphthoic anhydride, 1-naphthoic acid, and Bu3N were stirred at 160° for 4 h to give 85% (based on H3BO3) I (R1 = 1-naphthyl, R2 = R3 = R4 = Bu (II)). Then, a THF solution of the borate II was added dropwise to Me2CHOH solution of [Ph4P]Br at room temperature to give 85% K-free tetraphenylphosphonium tetrakis(1-naphthoyloxy)borate(1-).
IT **424822-39-9P**, 1-Benzyl-2-methylimidazolium tetrakis(trifluoroacetato)borate(1-) **424822-42-4P** **424822-44-6P**
RL: IMF (Industrial manufacture); SPN (Synthetic preparation); PREP (Preparation)
(preparation of tetrakis(acyloxy)borate from boric acid, anhydrides, and tertiary amines or ketimines and their substituted onium salts)
RN 424822-39-9 CAPLUS
CN Borate(1-), tetrakis(trifluoroacetato-κO)-, hydrogen, compd. with 2-methyl-1-(phenylmethyl)-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

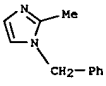
CM 1

CRN 36447-82-2
CMF C8 B F12 O8 . H
CCI CCS



CM 2

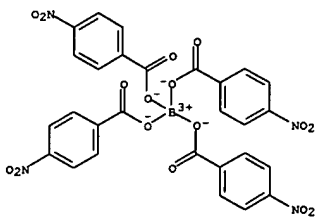
CRN 13750-62-4
CMF C11 H12 N2



RN 424822-42-4 CAPLUS
CN Borate(1-), tetrakis(4-nitrobenzoato-κO)-, hydrogen, compd. with 2,4,6-trimethylpyridine (1:1) (9CI) (CA INDEX NAME)

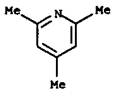
CM 1

CRN 347417-77-0
CMF C28 H16 B N4 O16 . H
CCI CCS



CM 2

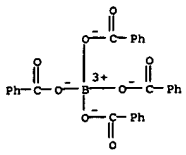
CRN 108-75-8
CMF C8 H11 N



RN 424822-44-6 CAPLUS
CN 1H-Imidazolium, 1,2,3-trimethyl-, tetrakis(benzoato-κO)borate(1-) (9CI) (CA INDEX NAME)

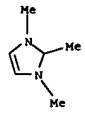
CM 1

CRN 182235-73-0
CMF C28 H20 B O8
CCI CCS



CM 2

CRN 65086-10-4
CMF C6 H11 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

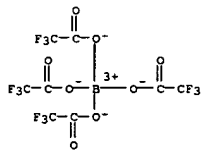
L12 ANSWER 25 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2002:358847 CAPLUS
DOCUMENT NUMBER: 136:369831
TITLE: Preparation of tetrakis(acyloxy) borates as curing accelerators for epoxy resins
INVENTOR(S): Nagata, Hiroshi
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 12 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2002138092	A2	20020514	JP 2000-330438	20001030
PRIORITY APPLN. INFO.:			JP 2000-330438	20001030

OTHER SOURCE(S): CASREACT 136:369831; MARPAT 136:369831
AB (R1CO2)4B-.N+HR2R3R4 (I; R1 = alkyl, aryl, aralkyl; R2-R4 = alkyl, aryl; R2R3 = :CR6R7; R6, R7 = alkyl, aryl; R2-R4, R6, R7 may form ring) are prepared by reaction of (R5O)3B (R5 = alkyl, aryl, aralkyl, acyl), (R1CO)2O (R1 = same as in I), R1CO2H (R1 = same as in I), and NR2R3R4 (R2-R4 = same as in I). E.g., tri-Et borate, 1-naphthoic anhydride, 1-naphthoic acid, and Bu3N were heated at 120° for 4 h to give 82% I (R1 = 1-naphthyl, R2-R4 = Bu).
IT **424822-39-9P 424822-42-4P 424822-44-6P**
RL: IMF (Industrial manufacture); SPN (Synthetic preparation); PREP (Preparation)
(preparation of tetrakis(acyloxy) borates as curing accelerators for epoxy resins)
RN 424822-39-9 CAPLUS
CN Borate(1-), tetrakis(trifluoroacetato-κO)-, hydrogen, compd. with 2-methyl-1-(phenylmethyl)-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

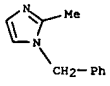
CM 1

CRN 36447-82-2
CMF C8 B F12 O8 . H
CCI CCS



CM 2

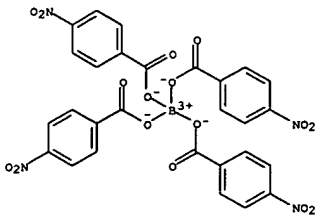
CRN 13750-62-4
CMF C11 H12 N2



RN 424822-42-4 CAPLUS
CN Borate(1-), tetrakis(4-nitrobenzoato-κO)-, hydrogen, compd. with 2,4,6-trimethylpyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

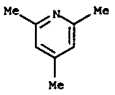
CRN 347417-77-0
CMF C28 H16 B N4 O16 . H
CCI CCS



● H+

CM 2

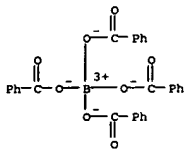
CRN 108-75-8
CMF C8 H11 N



RN 424822-44-6 CAPLUS
CN 1H-Imidazolium, 1,2,3-trimethyl-, tetrakis(benzoato-κO)borate(1-) (9CI) (CA INDEX NAME)

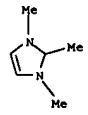
CM 1

CRN 182235-73-0
CMF C28 H20 B O8
CCI CCS



CM 2

CRN 65086-10-4
CMF C6 H11 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

L12 ANSWER 26 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

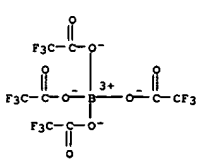
ACCESSION NUMBER: 2002:358846 CAPLUS
DOCUMENT NUMBER: 136:369830
TITLE: Preparation of tetrakis(acyloxy) borates as curing accelerators for epoxy resins
INVENTOR(S): Nagata, Hiroshi
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 11 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2002138091	A2	20020514	JP 2000-330140	20001030
PRIORITY APPLN. INFO.:			JP 2000-330140	20001030

OTHER SOURCE(S): CASREACT 136:369830; MARPAT 136:369830
AB (R1CO2)4B-.N+HR2R3R4 (I; R1 = alkyl, aryl, aralkyl; R2-R4 = alkyl, aryl; R2-R4 may form ring) are prepared by reaction of B2O3, (R1CO)2O (R1 = same as in I), R1CO2H (R1 = same as in I), and NR2R3R4 (R2-R4 = same as in I). E.g., B2O3, 1-naphthoic anhydride, 1-naphthoic acid, and Bu3N were heated at 160° for 4 h to give 85% I (R1 = 1-naphthyl, R2-R4 = Bu).
IT **424822-39-9P 424822-42-4P 424822-44-6P**
RL: IMF (Industrial manufacture); SPN (Synthetic preparation); PREP (Preparation)
(preparation of tetrakis(acyloxy) borates as curing accelerators for epoxy resins)
RN 424822-39-9 CAPLUS
CN Borate(1-), tetrakis(trifluoroacetato-κO)-, hydrogen, compd. with 2-methyl-1-(phenylmethyl)-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 36447-82-2
CMF C8 B F12 O8 . H
CCI CCS



● H+

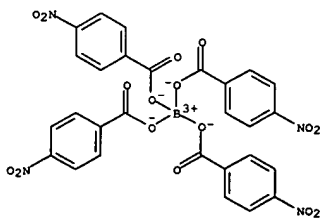
CM 2

CRN 13750-62-4
CMF C11 H12 N2

RN 424822-42-4 CAPLUS
CN Borate(1-), tetrakis(4-nitrobenzoato-κO)-, hydrogen, compd. with 2,4,6-trimethylpyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 347417-77-0
CMF C28 H16 B N4 O16 . H
CCI CCS

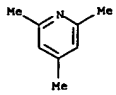


● H+

CM 2

CRN 108-75-8
CMF C8 H11 N

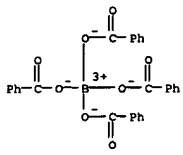
L12 ANSWER 26 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



RN 424822-44-6 CAPLUS
CN 1H-Imidazolium, 1,2,3-trimethyl-, tetrakis(benzoato-κO)borate(1-) (9CI) (CA INDEX NAME)

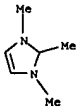
CM 1

CRN 182235-73-0
CMF C28 H20 B O8
CCI CCS



CM 2

CRN 65086-10-4
CMF C6 H11 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

L12 ANSWER 27 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2002:282748 CAPLUS
DOCUMENT NUMBER: 138:170125
TITLE: Hydrogen sulfate and tetrakis(hydrogen sulfato)borate ionic liquids: synthesis and catalytic application in highly Bronsted-acidic systems for Friedel-Crafts alkylation
AUTHOR(S): Wasserscheid, Peter; Sesing, Martin; Korth, Wolfgang
CORPORATE SOURCE: Institut fuer Technische Chemie und Makromolekulare Chemie der RWTH Aachen, Aachen, D-52074, Germany
SOURCE: Green Chemistry (2002), 4(2), 134-138
CODEN: GRCHFJ; ISSN: 1463-9262
PUBLISHER: Royal Society of Chemistry
DOCUMENT TYPE: Journal
LANGUAGE: English
OTHER SOURCE(S): CASREACT 138:170125

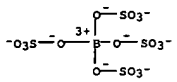
AB Hydrogen sulfate and tetrakis(hydrogen sulfato)borate ionic liqs. have been synthesized and characterized. These ionic liqs. are halogen-free, available from cheap raw materials and easy to prepare They have been used as additives to sulfuric acid in the Friedel-Crafts alkylation of benzene with 1-decene. The results clearly demonstrate an interplay of acidity and solubility effects caused by the ionic liquid additive. In some cases, low amts. of ionic liquid additive result in a dramatic improvement of product yield.

IT **497881-62-6P 497881-63-7P**
RL: CAT (Catalyst use); SPN (Synthetic preparation); PREP (Preparation); USES (Uses)
(preparation of imidazolium hydrogen sulfate and tetrakis(hydrogensulfato)borate ionic liqs. as catalysts for Friedel-Crafts alkylation)

RN 497881-62-6 CAPLUS
CN 1H-Imidazolium, 1-butyl-3-methyl-, hydrogen tetrakis[sulfato(2-)-κO]borate(5-) (1:4:1) (9CI) (CA INDEX NAME)

CM 1

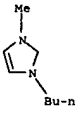
CRN 497881-61-5
CMF B O16 S4
CCI CCS



CM 2

CRN 80432-08-2
CMF C8 H15 N2

L12 ANSWER 27 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

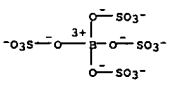


ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

RN 497881-63-7 CAPLUS
CN 1H-Imidazolium, 1-methyl-3-octyl-, hydrogen tetrakis[sulfato(2-)-κO]borate(5-) (1:4:1) (9CI) (CA INDEX NAME)

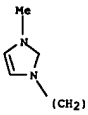
CM 1

CRN 497881-61-5
CMF B O16 S4
CCI CCS



CM 2

CRN 178631-03-3
CMF C12 H23 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE
REFERENCE COUNT: 24 THERE ARE 24 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 28 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2002:237754 CAPLUS
DOCUMENT NUMBER: 137:13532
TITLE: A polycyclic condensate of phenylboronic and boric acids with N-hydroxypiperidine
AUTHOR(S): Kliegel, Wolfgang; Drueckler, Klaus; Patrick, Brian O.; Rettig, Steven J.; Trotter, James
CORPORATE SOURCE: Technische Universitaet Braunschweig, Institut fuer Pharmazeutische Chemie, Braunschweig, 38106, Germany
SOURCE: Acta Crystallographica, Section E: Structure Reports Online (2002), E58(4), o393-o395
CODEN: ACSEBH; ISSN: 1600-5368
URL: http://journals.iucr.org/e/issues/2002/04/00/om60/81/om6081.pdf
PUBLISHER: International Union of Crystallography
DOCUMENT TYPE: Journal; (online computer file)
LANGUAGE: English

AB Crystals of 1-hydroxypiperidinium 16-hydroxy-6,6-pentamethylene-11,13,19-triphenyl-19-(1-piperidiniumoxy)-8,10,12,14,15,17,18,20-octaoxa-6-azonia-11,13,16-tribora-7,9,19-triboratatrispiro[5.0.1.5.3.3]eicosane Et2O hemisolvate are triclinic, space group P.hivin.1, with a 10.636(2), b 15.340(2), c 15.545(2) Å, α 105.467(2), β 107.949(4), γ 109.453(2)°; Z = 2, dc = 1.226; R = 0.043, Rw(F2) 0.102 for 8500 reflections. The compound contains a tetracyclic ring system, with two boroxine, a BOBON, and a piperidinium ring spiro-fused; the Et2O solvent mol. is disordered over an inversion center.

IT **433330-52-0P**, 1-Hydroxypiperidinium 16-hydroxy-6,6-pentamethylene-11,13,19-triphenyl-19-(1-piperidiniumoxy)-8,10,12,14,15,17,18,20-octaoxa-6-azonia-11,13,16-tribora-7,9,19-triboratatrispiro[5.0.1.5.3.3]eicosane compound with ethyl ether (2:1)
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(preparation and crystal structure of)

RN 433330-52-0 CAPLUS
CN Borate(2-), [μ-[1-(hydroxy-κO)piperidinato-κN]](1-(hydroxy-κO)piperidine][μ-[orthoborato(3-)-κO:κO']]di-μ-oxophenyl[(phenylboronic acid-κO) bimol. monoanhydridato(2-)]tri-, dihydrogen, compd. with 1-hydroxypiperidine and 1,1'-oxybis[ethane] (2:2:1) (9CI) (CA INDEX NAME)

CM 1

CRN 60-29-7
CMF C4 H10 O

CM 2

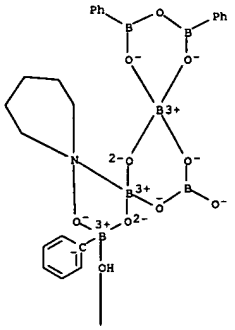
CRN 433330-51-9
CMF C28 H36 B6 N2 O10 . C5 H11 N O . 2 H

CM 3

CRN 433330-50-8
CMF C28 H36 B6 N2 O10 . 2 H

L12 ANSWER 28 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
CCI CCS

PAGE 1-A



PAGE 2-A

●2 H^+

CM 4

CRN 4801-58-5
CMF C5 H11 N O

L12 ANSWER 28 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE

FORMAT

L12 ANSWER 29 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2002:54507 CAPLUS
DOCUMENT NUMBER: 136:256313
TITLE: A Reduced Polyoxomolybdenum Borophosphate Anion Related to the Wells-Dawson Clusters
AUTHOR(S): Dumas, Eddy; Debiemme-Chouvy, Catherine; Sevov, Slavi C.
CORPORATE SOURCE: Department of Chemistry and Biochemistry, University of Notre Dame, Notre Dame, IN, 46556, USA
SOURCE: Journal of the American Chemical Society (2002), 124(6), 908-909
CODEN: JACSAT; ISSN: 0002-7863
PUBLISHER: American Chemical Society
DOCUMENT TYPE: Journal
LANGUAGE: English
OTHER SOURCE(S): CASREACT 136:256313

AB (C3N2H5)8[MoV5MoVI7O22(BO4)2(PO4)5(HPO4)3]·nH2O (n ≈ 4, C3N2H5 = imidazolium) is the 1st molybdenum borophosphate. It contains clusters of twelve molybdenum, eight phosphorus and two boron atoms, [Mo12O22(BO4)2(PO4)5(HPO4)3]8- similar to the Wells-Dawson clusters. Five molybdenum atoms are MoV while the others are MoVI, and the five electrons are delocalized. According to the temperature dependence of the magnetic susceptibility four of these electrons are paired.

IT **403824-78-2P**
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (preparation, crystal structure and magnetic susceptibility of polyoxomolybdenum borophosphate cluster)

RN 403824-78-2 CAPLUS
CN Molybdate(11-), diborateoctadeca-μ-oxododecaoxohexakis[μ3-{phosphato(3-)-κO:κO':κO'']]bis[μ7-{phosphato(3-)-κO:κO:κO':κO':κO'':κO'':κO''']]d odeca-, undecahydrogen, compd. with 1H-imidazole (1:8), tetrahydrate (9CI)
(CA INDEX NAME)

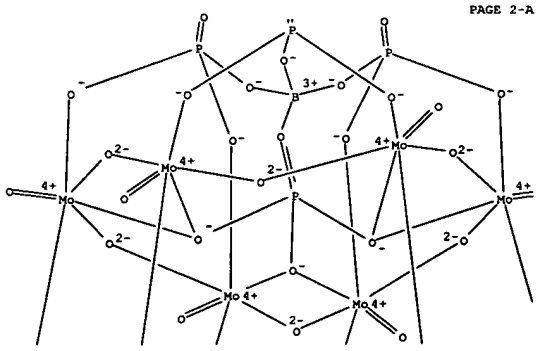
CM 1

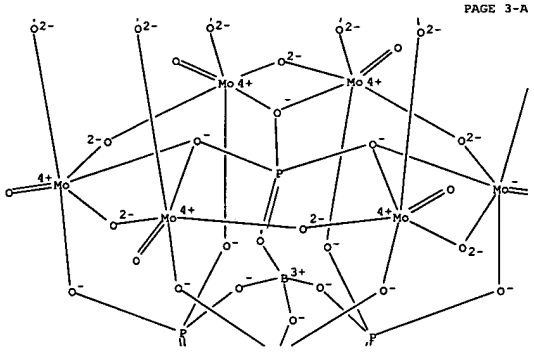
CRN 403824-77-1
CMF B2 Mo12 O62 P8 . 11 H
CCI CCS

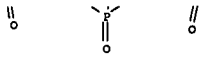
L12 ANSWER 29 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

PAGE 1-A

PAGE 2-A







●11 H^+

CM 2

CRN 288-32-4
CMF C3 H4 N2

REFERENCE COUNT: 31 THERE ARE 31 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE
FORMAT

L12 ANSWER 30 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2001:885417 CAPLUS
DOCUMENT NUMBER: 136:6139
TITLE: Ionic liquids
INVENTOR(S): Hilarius, Volker; Heider, Udo; Schmidt, Michael
PATENT ASSIGNEE(S): Merck Patent G.m.b.H., Germany
SOURCE: Eur. Pat. Appl., 27 pp.
CODEN: EPXXDW
DOCUMENT TYPE: Patent
LANGUAGE: German
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 1160249	A2	20011205	EP 2001-113237	20010530
EP 1160249	A3	20030917		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO				
DE 10026565	A1	20011206	DE 2000-10026565	20000530
BR 2001002084	A	20020319	BR 2001-2084	20010523
CA 2348966	AA	20011130	CA 2001-2348966	20010528
JP 2002187893	A2	20020705	JP 2001-158540	20010528
CN 1326936	A	20011219	CN 2001-119371	20010530
US 2002015883	A1	20020207	US 2001-866926	20010530
PRIORITY APPLN. INFO.:			DE 2000-10026565 A	20000530

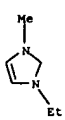
AB The preparation of title compds. is described. Thus, reaction of methylimidazole with chloroethane in MeCN gave ethylmethylimidazolium chloride which on basic hydrolysis gave 1-ethyl-3-methylimidazolium hydroxide. Reaction of 1-ethyl-3-methylimidazolium hydroxide with boric acid in presence of oxalic acid gave title compound, 1-ethyl-3-methylimidazoliumbis[1,2-oxalato(2-)-O,O']borate.

IT **376650-04-3P 376650-05-4P 376650-06-5P**
376650-07-6P
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation as ionic liquid)

RN 376650-04-3 CAPLUS
CN 1H-Imidazolium, 1-ethyl-3-methyl-, (T-4)-bis[1,2-benzenediolato(2-)-κO,κO']borate(1-) (9CI) (CA INDEX NAME)

CM 1

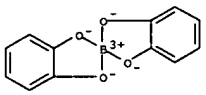
CRN 65039-03-4
CMF C6 H11 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

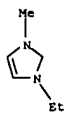
CRN 16986-25-7
CMF C12 H8 B O4
CCI CCS



RN 376650-05-4 CAPLUS
CN 1H-Imidazolium, 1-ethyl-3-methyl-, (T-4)-bis[2-(hydroxy-κO)benzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

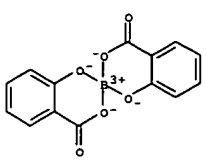
CRN 65039-03-4
CMF C6 H11 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

CRN 38403-08-6
CMF C14 H8 B O6
CCI CCS

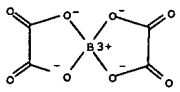


RN 376650-06-5 CAPLUS
CN 1H-Imidazolium, 1-ethyl-3-methyl-, (T-4)-bis[ethanedioato(2-)-κO1,κO2]borate(1-) (9CI) (CA INDEX NAME)

CM 1

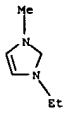
L12 ANSWER 30 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CRN 125579-65-9
CMF C4 B O8
CCI CCS



CM 2

CRN 65039-03-4
CMF C6 H11 N2

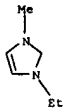


ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

RN 376650-07-6 CAPLUS
CN 1H-Imidazolium, 1-ethyl-3-methyl-, (T-4)-bis[2-(hydroxy-κO)propanoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

CRN 65039-03-4
CMF C6 H11 N2

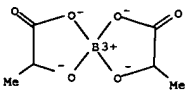


ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

CRN 31168-89-5
CMF C6 H8 B O6
CCI CCS

L12 ANSWER 30 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



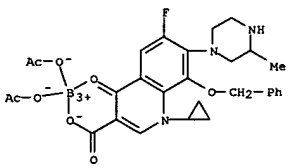
L12 ANSWER 31 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2001:859096 CAPLUS
DOCUMENT NUMBER: 137:201216
TITLE: New synthesis of Gatifloxacin
AUTHOR(S): Liu, Jiuyu; Tian, Zhiming; Guo, Huiyuan
CORPORATE SOURCE: Institute of Medicinal Biotechnology, Chinese Academy of medical Sciences and peking Union Medical College, Beijing, 100050, Peop. Rep. China
SOURCE: Zhongguo Yiyao Gongye Zazhi (2001), 32(10), 433-437
CODEN: ZYGZEA; ISSN: 1001-8255
PUBLISHER: Zhongguo Yiyao Gongye Zazhi Bianjibu
DOCUMENT TYPE: Journal
LANGUAGE: Chinese
OTHER SOURCE(S): CASREACT 137:201216

AB Gatifloxacin was synthesized from 3-hydroxy-2,4,5-trifluorobenzoic acid via 13 steps, with low overall yield. Ten new compds. were obtained, and their structures were characterized by 1HNMR and MS.

IT **452369-66-3P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(synthesis of Gatifloxacin)

RN 452369-66-3 CAPLUS
CN Boron, bis(acetato-κO)[1-cyclopropyl-6-fluoro-1,4-dihydro-7-(3-methyl-1-piperazinyl)-4-(oxo-κO)-8-(phenylmethoxy)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)



L12 ANSWER 32 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

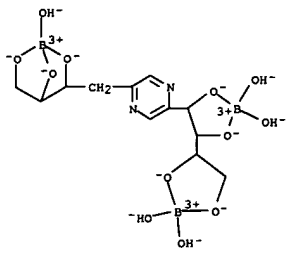
ACCESSION NUMBER: 2001:790434 CAPLUS
DOCUMENT NUMBER: 136:151372
TITLE: A clean conversion of D-glucosamine hydrochloride to a
pyrazine in the presence of phenylboronate or borate
AUTHOR(S): Rohovec, Jan; Kotek, Jan; Peters, Joop A.; Maschmeyer,
Thomas
CORPORATE SOURCE: Department of Applied Organic Chemistry and Catalysis,
Delft University of Technology, Delft, 2628 BL, Neth.
SOURCE: European Journal of Organic Chemistry (2001), (20), 3899-3901
CODEN: EJOCFK; ISSN: 1434-193X
PUBLISHER: Wiley-VCH Verlag GmbH
DOCUMENT TYPE: Journal
LANGUAGE: English
OTHER SOURCE(S): CASREACT 136:151372

AB D-Glucosamine was found to undergo a condensation to give 2-(arabino-tetrahydroxybutyl)-5-(erythro-2,3,4-trihydroxybutyl)-pyrazine as practically the sole product in the presence of phenylboronate or borate. The reaction proceeds in aqueous solns. at room temperature in 3 h in 58% isolated yield. In D2O solns., the incorporation of one deuterium into the methylene group of the trihydroxybutyl arm was found. The borate esters of the product were investigated by 11B and 13C NMR spectroscopy.

IT **394251-85-5P 394251-86-6P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(clean conversion of D-glucosamine hydrochloride to a pyrazine in the presence of phenylboronate or borate)

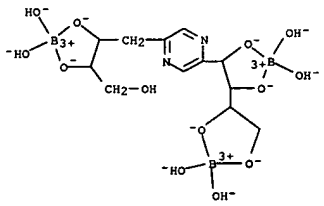
RN 394251-85-5 CAPLUS
CN Borate(3-), pentahydroxy[μ3-[(1R,2S,3R)-1-(5-[(2S,3R)-2,3,4-tri(hydroxy-κO)butyl]pyrazinyl]-1,2,3,4-butanetetrolato(7-)-κO1,κO2:κO3,κO4]]tri-, trisodium (9CI) (CA INDEX NAME)

L12 ANSWER 32 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



●3 Na+

RN 394251-86-6 CAPLUS
CN Borate(3-), hexahydroxy[μ3-[(1R,2S,3R)-1-[5-[(2S,3R)-2,3-di(hydroxy-κO)-4-hydroxybutyl]pyrazinyl]-1,2,3,4-butanetetrolato(6-)-κO1,κO2:κO3,κO4]]tri-, trisodium (9CI) (CA INDEX NAME)



●3 Na+

REFERENCE COUNT: 18 THERE ARE 18 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE
FORMAT

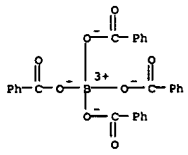
L12 ANSWER 33 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2001:704739 CAPLUS
DOCUMENT NUMBER: 135:273687
TITLE: Synthetic methods for tetrakis(acyloxy)borates and substituted onium tetrakis(acyloxy)borates
INVENTOR(S): Nagata, Hiroshi; Go, Yoshiyuki
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 8 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2001261684	A2	20010926	JP 2000-76942	20000317
PRIORITY APPLN. INFO.:			JP 2000-76942	20000317

OTHER SOURCE(S): MARPAT 135:273687
AB Boron trihalides react with carboxylic acid salts of amines or amidines to
prepare the title compds. Onium tetrakis(acyloxy)borates are useful for hardening accelerators for epoxy resins. Thus, triethylamine 1-naphthoic acid salt reacted with BBr3 to prepare triethylamine tetrakis(1-naphthoyloxy)borate.
IT **347417-74-7P 362588-75-8P**
RL: IMF (Industrial manufacture); PREP (Preparation)
(manufacture of tetrakis(acyloxy)borates and substituted onium tetrakis(acyloxy) borates)
RN 347417-74-7 CAPLUS
CN Borate(1-), tetrakis(benzoato-κO)-, hydrogen, compd. with 2-methyl-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 347417-73-6
CMF C28 H20 B O8 . H
CCI CCS



● H+

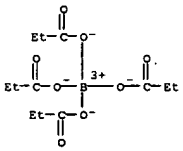
CM 2

L12 ANSWER 33 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
CRN 693-98-1
CMF C4 H6 N2

RN 362588-75-8 CAPLUS
CN 1H-Imidazolium, 1,2,3-trimethyl-, tetrakis(propanoato-κO)borate(1-) (9CI) (CA INDEX NAME)

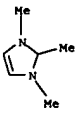
CM 1

CRN 362588-72-5
CMF C12 H20 B O8
CCI CCS



CM 2

CRN 65086-10-4
CMF C6 H11 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

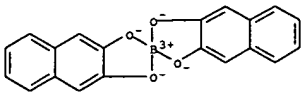
L12 ANSWER 34 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2001:684013 CAPLUS
DOCUMENT NUMBER: 135:227999
TITLE: Epoxy sealants with good curability and semiconductor devices using them
INVENTOR(S): Okubo, Akiko; Miyake, Sumiya
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 11 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2001253933	A2	20010918	JP 2000-65588	20000309
PRIORITY APPLN. INFO.:			JP 2000-65588	20000309

OTHER SOURCE(S): MARPAT 135:227999
AB The semiconductor sealants comprise (A) crystalline epoxy resins (m.p. 50-150°) having ≥2 epoxy groups in a mol., (B) compds. having ≥2 phenolic OH in a mol., (C) R1R2R3R4N+X- (I; R1-4 = H, monovalent aliphatic group, monovalent groups containing aromatic rings or
hetrorings; X = R5R6R7R8B; at least one of R5-8 = proton-released proton donor residues; rest of R5-8 = same as R1-4; elec. conductivity of the proton
donors ≤650 μS/cm for 2% aqueous solns. after a pressure cooker treatment), and (D) 200-2400 parts (based on 100 parts of A + B) inorg. fillers. Thus, a composition comprising 3,3',5,5'-tetramethylbiphenol diglycidyl ether polymer, HOC6H4[CH2-p-C6H4CH2C6H3(OH)]2.5H, I (R1-4 = Me,
X = Ph3BOQ, Q = 1-naphthalenyl; elec. conductivity of 1-naphthol 15 μS/cm),
fused spherical silica, and other additives showed spiral flow 93 cm, curing torque 90 kg-cm, spiral flow retention 90% after 1-wk storage, and good moisture resistance after cured.
IT **359765-78-9P 359767-63-8P**
RL: CAT (Catalyst use); IMF (Industrial manufacture); PREP (Preparation); USES (Uses)
(curing catalyst; epoxy sealants with good curability and storage stability for semiconductor devices)
RN 359765-78-9 CAPLUS
CN Pyrrolidinium, 1,1-dimethyl-, (T-4)-bis[2,3-naphthalenediolato(2-)-κO,κO']borate(1-) (9CI) (CA INDEX NAME)

CM 1

CRN 53992-98-6
CMF C20 H12 B O4
CCI CCS



CM 2

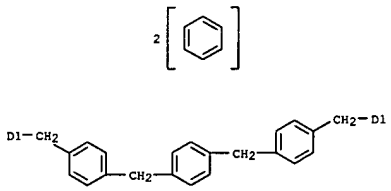
CRN 15312-12-6
CMF C6 H14 N

RN 359767-63-8 CAPLUS
CN Pyridinium, 1-methyl-, [2(or 5)-[[4-[[(hydroxy-κO)phenyl]methyl]phenyl]methyl]-5(or 2)-[[4-[(hydroxyphenyl)methyl]phenyl]methyl]phenolato]triphenoxyborate(1-) (9CI) (CA INDEX NAME)

CM 1

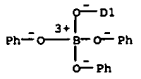
CRN 359767-62-7
CMF C52 H44 B O6
CCI CCS, IDS

PAGE 1-A



2 (D1—OH)

PAGE 2-A



CM 2

CRN 694-56-4
CMF C6 H8 N

L12 ANSWER 35 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2001:613298 CAPLUS
DOCUMENT NUMBER: 135:338424
TITLE: Acidity of heteropoly acids with various structures and compositions studied by IR spectroscopy of the pyridinium salts
AUTHOR(S): Maksimov, G. M.; Paukshtis, E. A.; Budneva, A. A.; Maksimovskaya, R. I.; Likholobov, V. A.
CORPORATE SOURCE: G. K. Boreskov Institute of Catalysis, Siberian Branch of the Russian Academy of Sciences, Novosibirsk, 630090, Russia
SOURCE: Russian Chemical Bulletin (Translation of Izvestiya Akademii Nauk, Seriya Khimicheskaya) (2001), 50(4), 587-590
CODEN: RCBUEY; ISSN: 1066-5285
PUBLISHER: Kluwer Academic/Consultants Bureau
DOCUMENT TYPE: Journal
LANGUAGE: English
AB The acidity on the proton affinity scale was determined by IR spectroscopy of the pyridinium salts for nineteen heteropoly acids of nine structural types (including two with the previously unknown structure) and one isopoly acid. All heteropoly acids exhibited a high acidity at the level of CF_3SO_3H and $HClO_4$. $H_3PW_{12}O_{40}$ was the strongest acid.
IT 368453-19-4
RL: PRP (Properties)
(acidity studied by IR spectroscopy)
RN 368453-19-4 CAPLUS
CN Tungstate(21-), henoctaconta-μ-oxononatriacontaoxotris[μ11-[tetrahydroxyborato(5-)-κO:κO:κO:κO':κO':.ka ppa.O':κO'':κO'':κO'':κO''':κO''']]nonatriacontaoxa-, heneicosahydrogen, compd. with pyridine (1:21) (9CI) (CA INDEX NAME)

CM 1

CRN 243445-97-8
CMF B3 O132 W39 . 21 H
CCI CCS

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

CM 2

CRN 110-86-1
CMF C5 H5 N

REFERENCE COUNT: 42 THERE ARE 42 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 36 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2001:545618 CAPLUS
DOCUMENT NUMBER: 135:139375
TITLE: Methods of preparation of nonaborate compositions
INVENTOR(S): Schubert, David M.
PATENT ASSIGNEE(S): U.S. Borax Inc., USA
SOURCE: PCT Int. Appl., 23 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

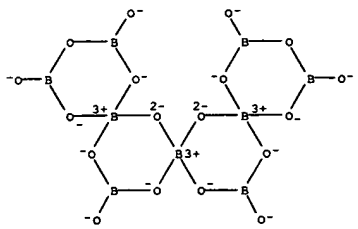
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001053201	A1	20010726	WO 2001-US1711	20010118
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW				
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
EP 1254075	A1	20021106	EP 2001-904921	20010118
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR				
JP 2003523964	T2	20030812	JP 2001-553215	20010118
TW 588015	B	20040521	TW 2001-90101327	20010119
US 2003109751	A1	20030612	US 2002-181628	20020719
US 6919036	B2	20050719		
PRIORITY APPLN. INFO.:			US 2000-177467P	P 20000121
			WO 2001-US1711	W 20010118

AB A family of borate compds. containing an isolated (finite) nonaborate anion with the structural formula $[B_9O_{12}(OH)_6]^{3-}$ is disclosed. Preferred amine nonaborate compds. have a resolved oxide formula of $A_2O.3B_2O_3.2H_2O$, where A is the monovalent cation of an amine salt, such as guanidinium and imidazolium. Also provided is a method for preparing these compds. by crystallization from an aqueous solution under mild conditions without the formation of significant amts. of byproducts. These compds. have potential application as precursors for the production of advanced boron nitride ceramic materials, and as flame retardants, corrosion inhibitors, and biocides.
IT 351428-05-2
RL: RCT (Reactant); RACT (Reactant or reagent)
(methods of preparation of nonaborate compns.)
RN 351428-05-2 CAPLUS
CN Borate(9-), bis[μ-[orthoborato(3-)-κO:κO']]di-μ-oxobis[μ-oxotetraoxodiborato(4-)]tri-, nonahydrogen, compd. with 1H-imidazole (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 273750-76-8
CMF B9 O18 . 9 H

L12 ANSWER 36 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
CCI CCS



●9 H+

CM 2

CRN 288-32-4
CMF C3 H4 N2

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE

FORMAT

L12 ANSWER 37 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2001:479794 CAPLUS
DOCUMENT NUMBER: 135:77683
TITLE: Manufacture of tetraacyl borates as crosslinking accelerators for epoxy resin for electric materials
INVENTOR(S): Nagata, Hiroshi; Miyake, Sumiya; Go, Yoshiyuki
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 13 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

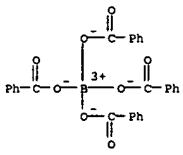
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2001181282	A2	20010703	JP 1999-371772	19991227
PRIORITY APPLN. INFO.:			JP 1999-371772	19991227

OTHER SOURCE(S): MARPAT 135:77683
AB The tetraacyl borates are manufactured by a reaction of (R1CO2)(R2CO2)(R3CO2)B with salts comprising NR4R5R6 and R7CO2H (R1-R3, R7 = alkyl, aryl, aralkyl; R4-R6 = alkyl, aryl; ≥2 of R4-R6 may form a ring). Thus, Et3N and a heptane solution of BBr3 were successively added into 1-naphthoic acid to give boron tri(1-naphthoate), which was heated with 1-naphthoic acid triethylamine salt to give 93% boron tetra(1-naphthoate) triethylamine salt.
IT **347417-74-7P**
RL: CAT (Catalyst use); IMF (Industrial manufacture); PREP (Preparation); USES (Uses)
(manufacture of tetraacyl borates as crosslinking accelerators for epoxy resins)
RN 347417-74-7 CAPLUS
CN Borate(1-), tetrakis(benzoato-κO)-, hydrogen, compd. with 2-methyl-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 347417-73-6
CMF C28 H20 B O8 . H
CCI CCS

L12 ANSWER 37 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



● H+

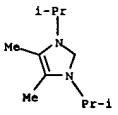
CM 2

CRN 693-98-1
CMF C4 H6 N2

L12 ANSWER 38 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2001:393042 CAPLUS
DOCUMENT NUMBER: 135:39092
TITLE: Crystal structure of 1,3-diisopropyl-4,5-dimethylimidazolium tetramethoxyborate, [C11H21N2][B(OCH3)4]
AUTHOR(S): Kuhn, N.; Steimann, M.; Weyers, G.
CORPORATE SOURCE: Universitat Tubingen, Institut fur Anorganische Chemie, Tubingen, D-72076, Germany
SOURCE: Zeitschrift fuer Kristallographie - New Crystal Structures (2001), 216(2), 315-317
CODEN: ZKNSFT; ISSN: 1433-7266
PUBLISHER: R. Oldenbourg Verlag
DOCUMENT TYPE: Journal
LANGUAGE: English
AB The title compound is triclinic, space group P.hivin.1, a 8.417(1), b 15.849(4), c 16.387(5) Å, α 61.37(1), β 85.73(2), γ 84.88(1)°, Z = 2, Rgt(F) = 0.057, wRref(F2) = 0.161, T = 293 K. Atomic coordinates are given. The shortest C-O distances is 3.007 Å.
IT **343781-24-8**
RL: PRP (Properties)
(crystal structure of)
RN 343781-24-8 CAPLUS
CN 1H-Imidazolium, 4,5-dimethyl-1,3-bis(1-methylethyl)-, tetramethoxyborate(1-) (9CI) (CA INDEX NAME)

CM 1

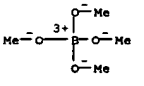
CRN 200803-10-7
CMF C11 H21 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

CRN 15841-16-4
CMF C4 H12 B O4
CCI CCS



REFERENCE COUNT: 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS

L12 ANSWER 38 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
RECORD. ALL CITATIONS AVAILABLE IN THE RE

FORMAT

L12 ANSWER 39 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2001:306245 CAPLUS
DOCUMENT NUMBER: 135:76743
TITLE: New Methods of Resolution and Purification of Racemic and Diastereomeric Amino Alcohol Derivatives Using Boric Acid and Chiral 1,1'-Bi-2-naphthol
AUTHOR(S): Periasamy, Mariappan; Kumar, Nangunoori Sampath; Sivakumar, Sangarappan; Rao, Vutukuri Dharma; Ramanathan, C. Ramaraj; Venkatraman, Lakshmanan
CORPORATE SOURCE: School of Chemistry, University of Hyderabad, Hyderabad, 500 046, India
SOURCE: Journal of Organic Chemistry (2001), 66(11), 3828-3833
CODEN: JOCEAH; ISSN: 0022-3263
PUBLISHER: American Chemical Society
DOCUMENT TYPE: Journal
LANGUAGE: English
OTHER SOURCE(S): CASREACT 135:76743

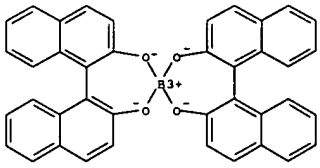
AB Resolution of racemic amino alc. derivs. is readily achieved to obtain enantiomerically enriched compds. using chiral 1,1'-bi-2-naphthol and boric acid in solvents such as MeCN, THF, and MeOH. Some of the intermediate ammonium borate complexes were also characterized by X-ray diffraction methods.

IT **347842-11-9P 347842-12-0P 347842-13-1P**
RL: PRP (Properties); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(resolution and purification of racemic and diastereomeric amino alc. derivs.
using boric acid and chiral 1,1'-bi-2-naphthol)

RN 347842-11-9 CAPLUS
CN Borate(1-), bis[[1,1'-binaphthalene]-2,2'-diolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with 1-(2-methoxycyclohexyl)piperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 347842-10-8
CMF C40 H24 B O4 . H
CCI CCS

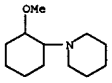


● H+

CM 2

L12 ANSWER 39 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

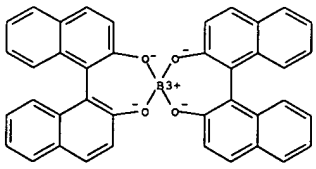
CRN 347841-84-3
CMF C12 H23 N O



RN 347842-12-0 CAPLUS
CN Borate(1-), bis[[1,1'-binaphthalene]-2,2'-diolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with (2R)-α,α-diphenyl-2-pyrrolidinemethanol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 347842-10-8
CMF C40 H24 B O4 . H
CCI CCS

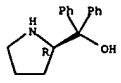


● H+

CM 2

CRN 22348-32-9
CMF C17 H19 N O

Absolute stereochemistry. Rotation (+).

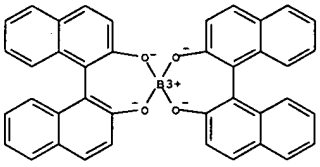


RN 347842-13-1 CAPLUS
CN Borate(1-), bis[[1,1'-binaphthalene]-2,2'-diolato(2-)-κO,κO']-

L12 ANSWER 39 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
, (T-4)-, hydrogen, compd. with acetonitrile and α-(1-methylethyl)-2-phenyl-1-pyrrolidineethanol (1:1:1) (9CI) (CA INDEX NAME)

CM 1

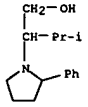
CRN 347842-10-8
CMF C40 H24 B O4 . H
CCI CCS



● H+

CM 2

CRN 347841-85-4
CMF C15 H23 N O



CM 3

CRN 75-05-8
CMF C2 H3 N

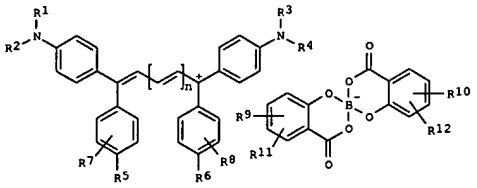
REFERENCE COUNT: 45 THERE ARE 45 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE

FORMAT

L12 ANSWER 40 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2000:762034 CAPLUS
DOCUMENT NUMBER: 133:323002
TITLE: Polymethine near infrared-absorbing dyes with high solubility for organic solvents
INVENTOR(S): Nakanishi, Isao; Saito, Nao
PATENT ASSIGNEE(S): Yamada Chemical Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 12 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2000302992	A2	20001031	JP 1999-152426	19990420
PRIORITY APPLN. INFO.:			JP 1999-152426	19990420

OTHER SOURCE(S): MARPAT 133:323002
GI



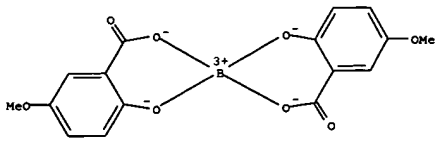
I

AB Title polymethine dyes are shown as I (R1-R4 = H, monovalent organic residue;
R1 and R2, or R3 and R4 may form ring; terminals of R1-R4 may form ring with C atoms in o-positions for N atoms bonded to aromatic groups; R5-R8 = H,
halo, monovalent organic residue; R9-R12 = H, OH, monovalent organic residue; n
= 0-3). The dyes are useful for near IR-absorbing filters, photothermal conversion materials, etc. Thus,
1,5-bis(p-dimethylaminophenyl)-1,5-bis(p-
tolyl)-2,4-pentadienol bis[3,5-bis(tert-butylsalicyl)]borate showed
λmax 814.5 nm and solubility 1.3% for acetone and 1.4% for CHCl3.
IT **303044-76-0P 303044-82-8P**
RL: IMF (Industrial manufacture); PRP (Properties); TEM (Technical or engineered material use); PREP (Preparation); USES (Uses)
(polymethine salicylborate near IR-absorbing dyes with high solubility for
organic solvents)
RN 303044-76-0 CAPLUS
CN Pyrrolidinium, 1-[4-[1,5,5-tris[4-(1-pyrrolidinyl)phenyl]-2,4-pentadienylidene]-2,5-cyclohexadien-1-ylidene]-, (T-4)-bis[2-(hydroxy-

L12 ANSWER 40 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
κO)-5-methoxybenzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

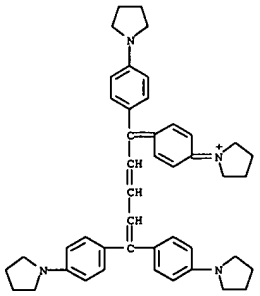
CM 1

CRN 303044-75-9
CMF C16 H12 B O8
CCI CCS



CM 2

CRN 162495-28-5
CMF C45 H51 N4

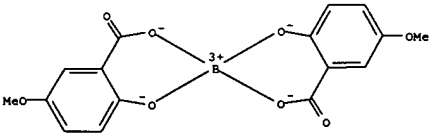


RN 303044-82-8 CAPLUS
CN Pyrrolidinium, 1-[4-[1,5-bis(4-methoxyphenyl)-5-[4-(1-pyrrolidinyl)phenyl]-2,4-pentadienylidene]-2,5-cyclohexadien-1-ylidene]-, (T-4)-bis[2-(hydroxy-
κO)-5-methoxybenzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

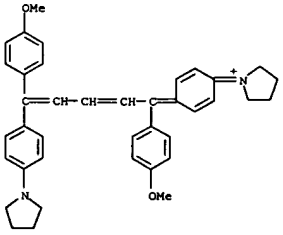
CRN 303044-75-9

L12 ANSWER 40 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
CMF C16 H12 B O8
CCI CCS



CM 2

CRN 303038-93-9
CMF C39 H41 N2 O2

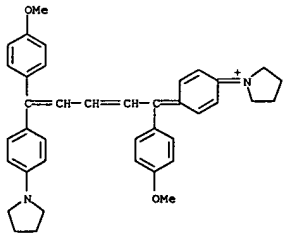


IT **303044-86-2**
RL: PRP (Properties); TEM (Technical or engineered material use); USES (Uses)
(polymethine salicylborate near IR-absorbing dyes with high solubility for
organic solvents)
RN 303044-86-2 CAPLUS
CN Pyrrolidinium, 1-[4-[1,5-bis(4-methoxyphenyl)-5-[4-(1-pyrrolidinyl)phenyl]-2,4-pentadienylidene]-2,5-cyclohexadien-1-ylidene]-, (T-4)-bis[2-(hydroxy-
κO)-3-methylbenzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

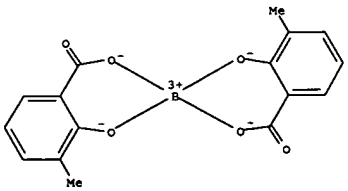
CRN 303038-93-9
CMF C39 H41 N2 O2

L12 ANSWER 40 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



CM 2

CRN 258875-08-0
CMF C16 H12 B O6
CCI CCS

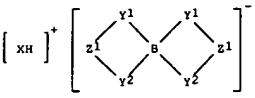


L12 ANSWER 41 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2000:657820 CAPLUS
DOCUMENT NUMBER: 133:253286
TITLE: Manufacture of latent crosslinking catalyst for thermosetting resins
INVENTOR(S): Go, Yoshiyuki; Miyake, Sumiya
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 8 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2000254513	A2	20000919	JP 1999-59915	19990308
PRIORITY APPLN. INFO.:			JP 1999-59915	19990308

OTHER SOURCE(S): MARPAT 133:253286
GI



I

AB The latent catalysts I are prepared by reacting tertiary amines or heterocyclic N-containing compds., with Y1Z1Y2 proton donors, and boric acid in solvents (Z1 = organic groups having substituents Y1, Y2; Y1, Y2 = groups that release proton and link and chelate with B atom; X = tertiary amines or heterocylic N-containing compds.). Stirring boric acid 12.4, salicylic acid 55.2, MeOH 276, and water 248 g for 30 min, adding a solution containing MeOH 378, water 378, and DBU 30.4 g and stirring for 1 h gave 80.3 g a white crystal onium borate of I with m.p. 128-130°.

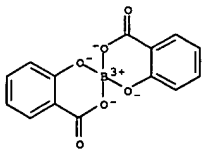
IT **233744-78-0P**
RL: CAT (Catalyst use); IMF (Industrial manufacture); PREP (Preparation); USES (Uses)
(manufacture of latent crosslinking catalyst for thermosetting resins)

RN 233744-78-0 CAPLUS
CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, (T-4)-, hydrogen, compd. with 2-methyl-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS

L12 ANSWER 41 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



CM 2

CRN 693-98-1
CMF C4 H6 N2

L12 ANSWER 42 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2000:626447 CAPLUS
DOCUMENT NUMBER: 134:36285
TITLE: Structural characterization of borate esters in which sodium acts as a support to the structural framework
AUTHOR(S): Bishop, Maximilienne; Bott, Simon G.; Barron, Andrew R.
CORPORATE SOURCE: Department of Chemistry, Rice University, Houston, TX, 77005, USA
SOURCE: Dalton (2000), (18), 3100-3105
CODEN: DALTFG
PUBLISHER: Royal Society of Chemistry
DOCUMENT TYPE: Journal
LANGUAGE: English
OTHER SOURCE(S): CASREACT 134:36285

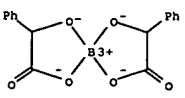
AB The borate ester derivs. of phenol, trans-cyclohexane-1,2-diol and mandelic acid were prepared from NaBH4 and structurally characterized by x-ray crystallog. The product from the reaction with phenol appears as a crystallog. disorder of [(THF)2Na(B(OPh)3H)]2 (1) and [(THF)2Na[B(OPh)(OH2)H]]2 (2). Both compds. are dimeric with bridging borate groups linking the Na cations. The reaction with trans-cyclohexane-1,2-diol in DMSO yields the infinite polymer, [(DMSO)Na(B(O2C6H10)2)]∞ (3), in which the Na cations link [B(O2C6H10)2]- anions. The unusual 5-coordinate geometry of the Na is completed by the coordination of a disordered DMSO mol. In a similar manner, mandelic acid reacts to form an infinite lattice [Na(py)2][B(O2CC(O)Ph)2] (4), in which each Na is coordinated to three [B(O2CC(O)Ph)2]- anions, one through two interactions involving the alkoxide and carboxylate groups of a chelate mandelic acid, and two interactions involving the carboxylate groups of adjacent anions. The role of the Group 1 cation in supporting the structural framework of the borate anions is discussed.

IT **312264-61-2P**
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(preparation and crystal structure of)

RN 312264-61-2 CAPLUS
CN Sodium(1+), bis(pyridine)-, (T-4)-bis[α-(hydroxy-κO)benzeneacetato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

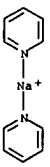
CRN 312264-60-1
CMF C16 H12 B O6
CCI CCS



CM 2

CRN 119327-39-8
CMF C10 H10 N2 Na
CCI CCS

L12 ANSWER 42 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



REFERENCE COUNT: 40 THERE ARE 40 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 43 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2000:428843 CAPLUS
DOCUMENT NUMBER: 133:159251
TITLE: Synthesis and Characterization of the Layered Vanadium Borophosphate (Imidazolium)3.8(H3O)1.2[(VO)4(BO)2(PO4)5]·0.3H2O
AUTHOR(S): Bontchev, Ranko P.; Do, Junghwan; Jacobson, Allan J.
CORPORATE SOURCE: Department of Chemistry, University of Houston, Houston, TX, 77204-5641, USA
SOURCE: Inorganic Chemistry (2000), 39(15), 3320-3324
CODEN: INOCAJ; ISSN: 0020-1669
PUBLISHER: American Chemical Society
DOCUMENT TYPE: Journal
LANGUAGE: English

AB The 1st layered V borophosphate (imidazolium)3.8(H3O)1.2[(VO)4(BO)2(PO4)5]·0.3H2O (1) was synthesized hydrothermally and characterized by chemical anal., IR and Raman spectroscopy, and thermogravimetric and magnetic measurements. The compound crystallizes in the monoclinic space group C2/c, a 9.4737(5) Å, b 22.1444(12) Å, c 17.2192(13) Å, β 105.936(1)°, Z = 4. The structure contains a novel borophosphate secondary building unit, [B2P5O22], in which two BP2O10 trimers are linked by an addnl. PO4 tetrahedron. These units are connected by V(IV)2O8 dimers and V(IV)O5 square pyramids to form layers. The space between the layers is filled by disordered imidazolium and hydronium cations and H2O mols. that form a complex network of H bonds. A model for the interlayer disorder is proposed.

IT **287470-82-0P**
RL: PRP (Properties); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation, crystal structure, antiferromagnetic exchange, vibrational spectra and thermal decomposition of)

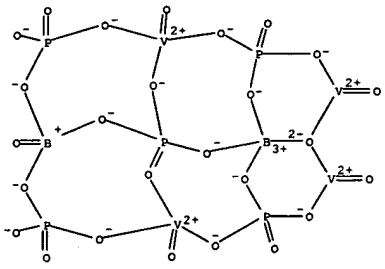
RN 287470-82-0 CAPLUS
CN Vanadate(5-), [μ4-[dioxo[μ-[phosphato(3-)-κO:κO']]tetrakis[phosphato(3-)-κO]diborato(13-)]]tetraoxotetra-, oxonium hydrogen, compd. with 1H-imidazole, hydrate (10:12:38:38:3) (9CI) (CA INDEX NAME)

CM 1

CRN 287470-81-9
CMF B2 O26 P5 V4 . 6/5 H3 O . 19/5 H
CCI CCS

L12 ANSWER 43 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

PAGE 1-A



PAGE 2-A

●19/5 H^+

●6/5 OH_3^+

CM 2

CRN 288-32-4
CMF C3 H4 N2

REFERENCE COUNT: 22 THERE ARE 22 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 44 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2000:278304 CAPLUS
DOCUMENT NUMBER: 133:37252
TITLE: Guanidinium and imidazolium borates containing the first examples of an isolated nonaborate oxoanion: [B9O12(OH)6]3-
AUTHOR(S): Schubert, David M.; Visi, Mandana Z.; Knobler, Carolyn B.
CORPORATE SOURCE: U.S. Borax Inc., Valencia, CA, 91355-1847, USA
SOURCE: Inorganic Chemistry (2000), 39(11), 2250-2251
CODEN: INOCAJ; ISSN: 0020-1669
PUBLISHER: American Chemical Society
DOCUMENT TYPE: Journal
LANGUAGE: English

AB The guanidinium and imidazolium borates [C(NH2)3]3[B9O12(OH)6] (I) and [C3H5N2]3[B9O12(OH)6] (II) were prepared and characterized by single crystal x-ray diffraction anal. These borates contain the 1st examples of the isolated nonaborate anion, [B9O12(OH)6]3-. These borates have resolved oxide formulas [C(NH2)3]2O·3B2O3·2H2O and [C3H5N2]2O·3B2O3·2H2O. The [B9O12(OH)6]3- anion found in each consists of four B3O3 rings sharing three tetrahedral B centers in a linear arrangement. The remaining six B centers are trigonal with one attached hydroxyl group. The two inner B3O3 rings each contain one trigonal and two tetrahedral B's, and the two outer rings each contain one tetrahedral and two trigonal B's.

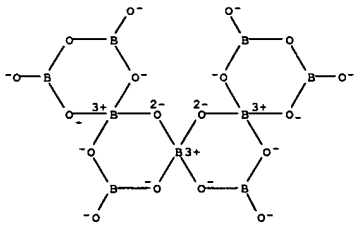
IT **273750-78-0P**
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(preparation and crystal structure of)

RN 273750-78-0 CAPLUS
CN Borate(9-), bis[μ-[orthoborato(3-)-κO:κO']]di-μ-oxobis[μ-oxotetraoxodiborato(4-)]tri-, nonahydrogen, compd. with 1H-imidazole (1:9) (9CI) (CA INDEX NAME)

CM 1

CRN 273750-76-8
CMF B9 O18 . 9 H
CCI CCS

L12 ANSWER 44 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



●9 H^+

CM 2

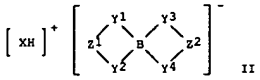
CRN 288-32-4
CMF C3 H4 N2

REFERENCE COUNT: 8 THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 45 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2000:37940 CAPLUS
DOCUMENT NUMBER: 132:79372
TITLE: Thermosetting novolak resin compositions with good curability, and molding materials therefrom
INVENTOR(S): Oka, Wataru; Orihara, Tamotsu
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 6 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2000017145	A2	20000118	JP 1998-184487	19980630
PRIORITY APPLN. INFO.:			JP 1998-184487	19980630

OTHER SOURCE(S): MARPAT 132:79372
GI



II

AB The compns. contain novolak resins, hexamethylenetetramine (I), and onium borates II (X = N-containing heterocyclic compound; Z1, Z2 = aromatic or alicyclic group; Y1-Y4 = proton donor group residue). Thus, novolak resin 100, I 16, II (X = DBU, Y1Z1Y2 = Y3Z2Y4 = O-o-C6H4CO2) 8 parts, and fillers are mixed and transfer-molded to give a test piece showing Barcol hardness 68 and bending strength 102 and 63 MPa, at room temperature and 120°, resp.

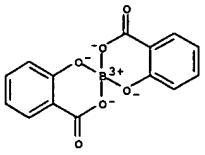
IT **233744-78-0**
RL: CAT (Catalyst use); USES (Uses)
(curing accelerator; thermosetting novolak resin compns. with good curability)

RN 233744-78-0 CAPLUS
CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, (T-4)-, hydrogen, compd. with 2-methyl-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS

L12 ANSWER 45 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



● H+

CM 2

CRN 693-98-1
CMF C4 H6 N2

L12 ANSWER 46 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1999:536707 CAPLUS
DOCUMENT NUMBER: 131:299232
TITLE: Resolution of trans-(±)-2-(pyrrolidinyl)cyclohexanol and its methyl ether using boric acid and chiral 1,1'-bi-2-naphthol
AUTHOR(S): Periasamy, Mariappan; Ramanathan, C. Ramaraj; Kumar, Nangunoori Sampath
CORPORATE SOURCE: School of Chemistry, University of Hyderabad, Hyderabad, 500 046, India
SOURCE: Tetrahedron: Asymmetry (1999), 10(12), 2307-2310
CODEN: TASYE3; ISSN: 0957-4166
PUBLISHER: Elsevier Science Ltd.
DOCUMENT TYPE: Journal
LANGUAGE: English

AB Trans-(±)-2-(pyrrolidinyl)cyclohexanol and its Me ether were resolved using chiral 1,1'-bi-2-naphthol and boric acid in THF or CH3CN. X-ray structural anal. was carried out for a diastereomeric complex obtained using the Me ether.

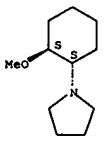
IT **247114-34-7P**
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(resolution of trans-(pyrrolidinyl)cyclohexanol and its Me ether using boric acid and chiral binaphthol)

RN 247114-34-7 CAPLUS
CN Borate(1-), bis[(1R)-[1,1'-binaphthalene]-2,2'-diolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with acetonitrile and 1-[(1S,2S)-2-methoxycyclohexyl]pyrrolidine (1:3:1) (9CI) (CA INDEX NAME)

CM 1

CRN 247024-88-0
CMF C11 H21 N O

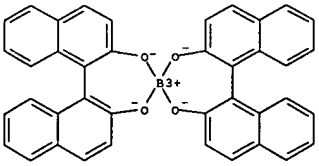
Absolute stereochemistry. Rotation (+).



CM 2

CRN 161651-51-0
CMF C40 H24 B O4 . H
CCI CCS

L12 ANSWER 46 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



● H+

CM 3

CRN 75-05-8
CMF C2 H3 N

REFERENCE COUNT: 15 THERE ARE 15 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 47 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1999:481505 CAPLUS
DOCUMENT NUMBER: 131:130718
TITLE: Epoxy resin compositions for prepregs and laminated circuit boards
INVENTOR(S): Go, Yoshiyuki; Miyake, Sumiya; Nagata, Hiroshi; Okubo, Akiko; Kobayashi, Minoru
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 8 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 11209583	A2	19990803	JP 1998-11829	19980123
PRIORITY APPLN. INFO.:			JP 1998-11829	19980123

OTHER SOURCE(S): MARPAT 131:130718

AB Title epoxy resin compns., which is stable at room temperature and fat curing upon heating, comprise an epoxy resin, a polyamine curing agent, and a curing accelerator of an onium borate. Thus a bisphenol A-based epoxy resin 100 parts, diaminodiphenylmethane 0.2 parts, dicyandiamide 0.3 parts, and DBU salt of 2-hydroxybenzoic acid boron complex 2.5 parts were mixed to give an epoxy composition of this invention. Two pieces of glass cloths were impregnated with the above composition to two prepregs which were laminated with an printed circuit board by vacuum compression while heating to give a laminated board.

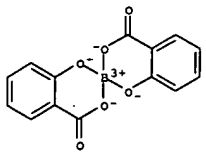
IT **233744-78-0 233744-79-1 233744-80-4**
RL: CAT (Catalyst use); USES (Uses)
(epoxy resin compns. for prepregs and laminated circuit boards)

RN 233744-78-0 CAPLUS
CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, (T-4)-, hydrogen, compd. with 2-methyl-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS

L12 ANSWER 47 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



● H+

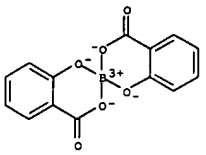
CM 2

CRN 693-98-1
CMF C4 H6 N2

RN 233744-79-1 CAPLUS
CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, (T-4)-, hydrogen, compd. with 4-methyl-2-phenyl-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS

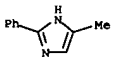


● H+

L12 ANSWER 47 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CM 2

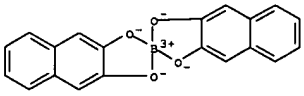
CRN 827-43-0
CMF C10 H10 N2



RN 233744-80-4 CAPLUS
CN Borate(1-), bis[2,3-naphthalenediolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with 2-methyl-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 47422-29-7
CMF C20 H12 B O4 . H
CCI CCS



● H+

CM 2

CRN 693-98-1
CMF C4 H6 N2

L12 ANSWER 48 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1999:470222 CAPLUS
DOCUMENT NUMBER: 131:208123
TITLE: Templated synthesis of vanadium borophosphate cluster anions
AUTHOR(S): Bontchev, Ranko P.; Do, Junghwan; Jacobson, Allan J.
CORPORATE SOURCE: Department of Chemistry, University of Houston, Houston, TX, 77204-5641, USA
SOURCE: Angewandte Chemie, International Edition (1999), 38(13/14), 1937-1940
CODEN: ACIEF5; ISSN: 1433-7851
PUBLISHER: Wiley-VCH Verlag GmbH
DOCUMENT TYPE: Journal
LANGUAGE: English

AB Hydrothermal reactions of V2O3, H3BO3 and H3PO4 at 180° for 3 days gave crystals of (NMe4)6[(VO)2BP2O10]4·nH2O (1; n = 2,6,14), Na14[Na ⊃ [(VO)2BP2O10)5]·nH2O (2) and A17[A ⊃ [(VO)2BP2O10)6]·nH2O (3; A = NH4, K, Rb, Cs). The crystal structures of 1 (n = 6), 2 and 3 (A = NH4) and the magnetic susceptibility of 3 (indicating antiferromagnetic interactions)are reported. Thermogravimetric anal. of 1-3 indicate structure collapse upon loss of H2O with eventual formation of glassy matrixes composed of VO2, P5O5, B2O3 and M2O (M = Na, K, Rb, Cs).

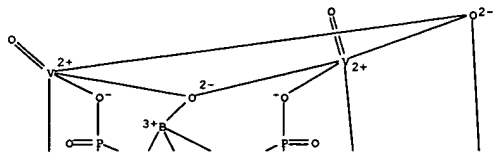
IT **241125-89-3P**
RL: PEP (Physical, engineering or chemical process); PRP (Properties); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); PROC (Process); RACT (Reactant or reagent)
(preparation and crystal and mol. structure and thermal decomposition)

RN 241125-89-3 CAPLUS
CN Vanadate(12-), tetraborateocta-μ3-oxooctaoxooctakis[μ3-[phosphato(3-)-κO:κO':κO'']]octa-, dodecahydrogen, compd. with piperazine (1:6), hexahydrate (9CI) (CA INDEX NAME)

CM 1

CRN 241125-88-2
CMF B4 O48 P8 V8 . 12 H
CCI CCS

PAGE 1-A



* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

PAGE 4-A

●12 H^+

CM 2

CRN 110-85-0
CMF C4 H10 N2

IT **241125-95-1P 241125-96-2P**
RL: PEP (Physical, engineering or chemical process); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); PROC (Process); RACT

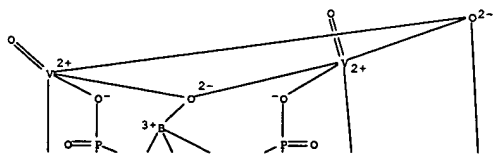
CRN 110-85-0
CMF C4 H10 N2

RN 241125-96-2 CAPLUS
CN Vanadate(12-), tetraborateocta-μ3-oxooctaoxooctakis[μ3-{phosphato(3-)-κO:κO':κO''}]octa-, dodecahydrogen, compd. with piperazine (1:6), tetradecahydrate (9CI) (CA INDEX NAME)

CM 1

CRN 241125-88-2
CMF B4 O48 P8 V8 . 12 H
CCI CCS

PAGE 1-A



* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

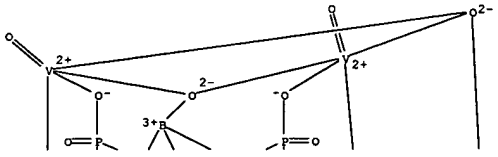
* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

(Reactant or reagent)
(prepn. and thermal decompn.)
RN 241125-95-1 CAPLUS
CN Vanadate(12-), tetraborateocta-μ3-oxooctaoxooctakis[μ3-{phosphato(3-)-κO:κO':κO''}]octa-, dodecahydrogen, compd. with piperazine (1:6), dihydrate (9CI) (CA INDEX NAME)

CM 1

CRN 241125-88-2
CMF B4 O48 P8 V8 . 12 H
CCI CCS

PAGE 1-A



* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

PAGE 4-A

●12 H^+

CM 2

PAGE 4-A

●12 H^+

CM 2

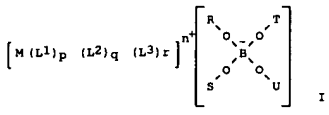
CRN 110-85-0
CMF C4 H10 N2

REFERENCE COUNT: 43 THERE ARE 43 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 49 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1999:463430 CAPLUS
DOCUMENT NUMBER: 131:151739
TITLE: Metal ion source for improving image-fixability
INVENTOR(S): Okubo, Kimihiko; Asatake, Atsushi
PATENT ASSIGNEE(S): Konica Co., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 10 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 11199790	A2	19990727	JP 1998-12169	19980106
PRIORITY APPLN. INFO.:			JP 1998-12169	19980106

OTHER SOURCE(S): MARPAT 131:151739
GI



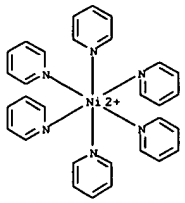
I

AB The metal ion source for improving dye-fixability is represented by a general formula I (M = metal ion; L1-3 = ligand; p = 0-3; q = 0-2; r = 0, 1; n = 1-3; R, S, T, U = H, substituent; 2 of R, S, T, U may form ring). The metal ion source shows improved storage stability.
IT **235782-66-8**
RL: MOA (Modifier or additive use); USES (Uses)
(metal ion source with improved storage stability for improving image-fixability)
RN 235782-66-8 CAPLUS
CN Nickel(2+), hexakis(pyridine)-, (OC-6-11)-, bis[(T-4)-bis[1,2-benzenediolato(2-)-κO,κO']borate(1-)] (9CI) (CA INDEX NAME)

CM 1

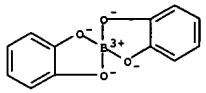
CRN 20037-72-3
CMF C30 H30 N6 Ni
CCI CCS

L12 ANSWER 49 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



CM 2

CRN 16986-25-7
CMF C12 H8 B O4
CCI CCS



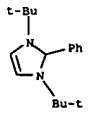
L12 ANSWER 50 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1999:407164 CAPLUS
DOCUMENT NUMBER: 131:103041
TITLE: Storage-stable epoxy resin compositions containing ammonium borates as latent crosslinking accelerators
INVENTOR(S): Miyake, Sumiya; Go, Yoshiyuki; Nagata, Hiroshi; Okubo, Akiko; Kobayashi, Minoru
PATENT ASSIGNEE(S): Sumitomo Bakelite Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 8 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 11171981	A2	19990629	JP 1997-340126	19971210
JP 3690710	B2	20050831		
PRIORITY APPLN. INFO.:			JP 1997-340126	19971210

OTHER SOURCE(S): MARPAT 131:103041
GI For diagram(s), see printed CA Issue.
AB Title compns., useful for elec. an electronic devices, etc., contain hardeners and X+ BY1Y2Y3Y4- [X+ = (substituted) ammonium; ≥1 of Y1-Y4 = H+-donating group residue after releasing 1 H+, the rest of Y1-Y4 = aromatic, heterocyclic, or aliphatic group] or I (Y9-Y10 are same as Y1-Y4; Y11-Y12 = H+-donating group residue after release of H+). Thus, o-cresol novolak epoxy resin (EOCN 102065) 67, phenol novolak 33, pulverized fused silica 300, carnauba wax 2, and Ph4N+ (BzO)4B- 3.1 part was mixed and roll-kneaded at 90° for 5 min to give title composition having initial spiral flow 83 cm and 79 cm after 3-day storage at 40°.
IT **229316-59-0 229316-71-6**
RL: CAT (Catalyst use); USES (Uses)
(rapidly curable storage-stable epoxy resin compns. containing ammonium borates as latent crosslinking accelerators)
RN 229316-59-0 CAPLUS
CN 1H-Imidazolium, 1,3-bis(1,1-dimethylethyl)-2-phenyl-, (T-4)-bis[1,2-benzenediolato(2-)-κO,κO']borate(1-) (9CI) (CA INDEX NAME)

CM 1

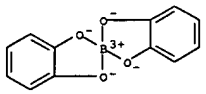
CRN 229316-58-9
CMF C17 H25 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

L12 ANSWER 50 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
CM 2

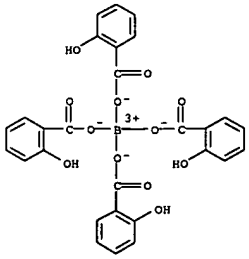
CRN 16986-25-7
CMF C12 H8 B O4
CCI CCS



RN 229316-71-6 CAPLUS
CN Pyridinium, 1-ethyl-, tetrakis(2-hydroxybenzoato-κO)borate(1-) (9CI) (CA INDEX NAME)

CM 1

CRN 229316-70-5
CMF C28 H20 B O12
CCI CCS



CM 2

CRN 15302-96-2
CMF C7 H10 N

ACCESSION NUMBER: 1999:163192 CAPLUS
DOCUMENT NUMBER: 130:245332
TITLE: Electrolytic capacitor driving electrolytic solution containing borodisalicylate
INVENTOR(S): Uramoto, Masahide; Nakano, Minoru; Sano, Mikio; Takahashi, Isao; Fukuda, Mitsuru
PATENT ASSIGNEE(S): Toyama Yakuhin Kogyo K. K., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 5 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 11067604	A2	19990309	JP 1997-225608	19970808
PRIORITY APPLN. INFO.:			JP 1997-225608	19970808

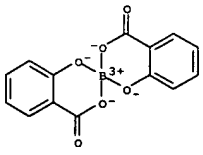
OTHER SOURCE(S): MARPAT 130:245332
GI

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The solution contains a borodisalicylic acid amine salt, as an electrolyte, comprising (A) a borodisalicylic acid backbone I and (2) an amine compound backbone NR1R2R3, II, or III (R1-5 = H, C1-5 alkyl; C1-7 alkylamino, cyclic group). The solution shows low resistivity, heat stability, and less odor.
IT **221332-52-1 221332-56-5**
RL: DEV (Device component use); MOA (Modifier or additive use); USES (Uses)
(electrolytic capacitor driving solution containing borodisalicylic acid amine derivative)
RN 221332-52-1 CAPLUS
CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, (T-4)-, hydrogen, compd. with 1,2-dimethyl-1H-imidazole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS



● H+

CM 2

CRN 1739-84-0
CMF C5 H8 N2

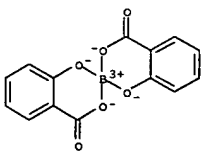
RN 221332-56-5 CAPLUS
CN Pyrimidinium, 1,2-dimethyl-, (T-4)-bis[2-(hydroxy-κO)benzoato(2-)-κO]borate(1-) (9CI) (CA INDEX NAME)

CM 1

CRN 221332-55-4
CMF C6 H9 N2

CM 2

CRN 38403-08-6
CMF C14 H8 B O6
CCI CCS



L12 ANSWER 52 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1999:139922 CAPLUS
DOCUMENT NUMBER: 130:189140
TITLE: Electrochromic system with coupled red-ox system and special anions
INVENTOR(S): Berneth, Horst; Kostromine, Serguei
PATENT ASSIGNEE(S): Bayer A.-G., Germany
SOURCE: PCT Int. Appl., 93 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: German
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9909111	A1	19990225	WO 1998-EP4909	19980806
W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
DE 19735733	A1	19990225	DE 1997-19735733	19970818
CA 2300131	AA	19990225	CA 1998-2300131	19980806
AU 9893393	A1	19990308	AU 1998-93393	19980806
AU 733812	B2	20010524		
EP 1015524	A1	20000705	EP 1998-946284	19980806
R: AT, DE, ES, FR, GB, IT, NL, SE, IE				
JP 2001515120	T2	20010918	JP 2000-509781	19980806
US 6417951	B1	20020709	US 2000-485758	20000215
US 2002197486	A1	20021226	US 2002-150335	20020517
US 6767481	B2	20040727		

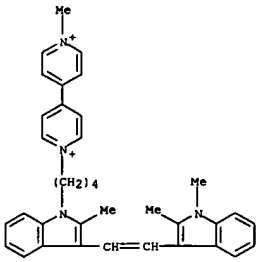
PRIORITY APPLN. INFO.: DE 1997-19735733 A 19970818
WO 1998-EP4909 W 19980806
US 2000-485758 A3 20000215

OTHER SOURCE(S): MARPAT 130:189140
AB Electrochromic systems are described which comprise ≥1 reducible substance and ≥1 oxidizable substance interconnected via a covalent binding link. Preferably the anions have mol. wts. >200 g/mol (most preferably >250 g/mol) and/or are cage-type anions. Electrochromic fluids comprising ≥1 of the compds. in ≥1 inert solvent are also described, as are electrochromic devices (e.g., solar cells, windows, mirrors, sun roofs, and displays) using the fluids.
IT **220624-16-8P**
RL: DEV (Device component use); PRP (Properties); SPN (Synthetic preparation); PREP (Preparation); USES (Uses)
(electrochromic materials with coupled red-ox systems and heavy and/or cage-type anions and devices using them)
RN 220624-16-8 CAPLUS
CN 4,4'-Bipyridinium, 1-[4-[3-[2-(1,2-dimethyl-1H-indol-3-yl)ethenyl]-2-methyl-1H-indol-1-yl]butyl]-1'-methyl-, bis[tetrabutoxyborate(1-)] (9CI) (CA INDEX NAME)

CM 1

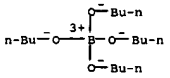
L12 ANSWER 52 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CRN 195328-26-8
CMF C36 H38 N4



CM 2

CRN 103624-08-4
CMF C16 H36 B O4
CCI CCS

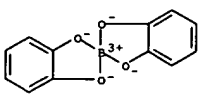


REFERENCE COUNT: 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 53 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1999:82369 CAPLUS
DOCUMENT NUMBER: 130:117629
TITLE: Salts of the bis(catecholato)borate anion with organic cations
AUTHOR(S): Clegg, William; Scott, Andrew J.; Lawlor, Fiona J.; Norman, Nicholas C.; Marder, Todd B.; Dai, Chaoyang; Nguyen, Paul
CORPORATE SOURCE: Department of Chemistry, University of Newcastle upon Tyne, Newcastle upon Tyne, NE1 7RU, UK
SOURCE: Acta Crystallographica, Section C: Crystal Structure Communications (1998), C54(12), 1875-1880
CODEN: ACSCEE; ISSN: 0108-2701
PUBLISHER: Munksgaard International Publishers Ltd.
DOCUMENT TYPE: Journal
LANGUAGE: English
AB In six salts with organic N and P cations [2-methylpyridinium bis(pyrocatecholato-O,O')borate, C6H8N+·C12H8BO4-, (1); 4-methylpyridinium bis(pyrocatecholato-O,O')borate, C6H8N+·C12H8BO4- (2) and (3) (two polymorphs); 1,10-phenanthrolinium bis(pyrocatecholato-O,O')borate, C12H9N2+·C12H8BO4-, (4), and its CH2Cl2 solvate, C12H9N2+·C12H8BO4-·CH2Cl2, (5); and trimethylphosphonium bis(pyrocatecholato-O,O')borate, C3H10P+·C12H8BO4-, (6)], the bis(catecholato)borate anion has approx. D2d (42m) symmetry, with the central spiro-B atom distorted from regular tetrahedral coordination geometry by reduction of the two intra-ring O-B-O bond angles. The two chelate rings show small deviations from planarity by folding about the O···O axis. Ion pairs are formed by N-H···O H bonding in all five salts with N-based cations, but there is no H bonding in the phosphonium salt; the H bonding leads to slight elongation of the B-O bond involved. Crystallog. data are given.
IT **30776-62-6**, 2-Methylpyridinium bis(pyrocatecholato-O,O')borate
RL: PRP (Properties)
(crystal structure of)
RN 30776-62-6 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with 2-methylpyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H+

L12 ANSWER 53 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

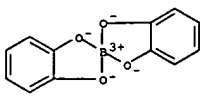
CM 2

CRN 109-06-8
CMF C6 H7 N

IT **219702-48-4**, 4-Methylpyridinium bis(pyrocatecholato-O,O')borate
RL: PRP (Properties)
(crystal structure of polymorphs of)
RN 219702-48-4 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with 4-methylpyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H+

CM 2

CRN 108-89-4
CMF C6 H7 N

REFERENCE COUNT: 15 THERE ARE 15 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

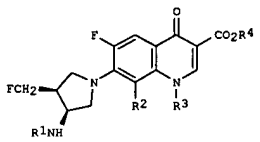
L12 ANSWER 53 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

L12 ANSWER 54 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1999:27822 CAPLUS
DOCUMENT NUMBER: 130:81423
TITLE: Preparation of cis-substituted fluoromethylpyrrolidine derivatives of 1,4-dihydro-4-oxoquinoline-3-carboxylic acid as antibacterial agents
INVENTOR(S): Takemura, Makoto; Takahashi, Hisashi; Ohki, Hitoshi; Kimura, Kenichi; Miyauchi, Rie; Takeda, Toshiyuki
PATENT ASSIGNEE(S): Daiichi Pharmaceutical Co., Ltd., Japan
SOURCE: PCT Int. Appl., 51 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9858923	A1	19981230	WO 1998-JP2787	19980623
W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, GM, GW, HU, ID, IL, IS, JP, KE, KG, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
AU 9880387	A1	19990104	AU 1998-80387	19980623
ZA 9805466	A	19990120	ZA 1998-5466	19980623
EP 995744	A1	20000426	EP 1998-928627	19980623
EP 995744	B1	20030212		
R: BE, CH, DE, FR, GB, IT, LI, NL, SE				
TW 382625	B	20000221	TW 1998-87110150	19980624
NO 9906390	A	20000224	NO 1999-6390	19991222
US 2002072608	A1	20020613	US 1999-446696	19991223
US 6656952	B2	20031202		
PRIORITY APPLN. INFO.:			JP 1997-166438 A	19970624
			JP 1998-54700 A	19980306
			WO 1998-JP2787 W	19980623

OTHER SOURCE(S): MARPAT 130:81423
GI



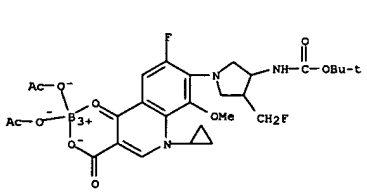
I

L12 ANSWER 54 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

AB Quinolone derivs. represented by general formula (I; wherein R1 represents a hydrogen atom or an optionally substituted alkyl group having 1 to 6 carbon atoms; R2 represents a halogenomethoxyl or alkoxyl group; R3 represents an alkyl, alkenyl, halogenoalkyl, cyclic alkyl, heteroaryl, alkoxyl, or alkylamino group; R4 represents a hydrogen atom or Ph, acetoxymethyl, pivaloyloxymethyl, ethoxycarbonyl, choline, dimethylaminoethyl, 5-indanyl, phthalidinyl, 5-alkyl-2-oxo-1,3-dioxol-4-ylmethyl, 3-acetoxy-2-oxobutyl, alkyl, alkoxymethyl, or phenylalkyl group), salts thereof, and hydrates thereof of both, having potent antimicrobial activities against various bacteria including resistant strains and being highly safe, are prepared Thus, 327 mg 3-(S)-tert-butoxycarbonylamino-4-(S)-fluoromethylpyrrolidine and 400 μL Et3N were added to a solution of 345 mg 1-cyclopropyl-6,7-difluoro-1,4-dihydro-8-methoxy-4-oxoquinoline-3-carboxylic acid-BF2 chelate in 2 mL DMSO, followed by deprotection with concentrated HCl to give 47% I (R1 = R4 = H, R2 = OMe, R3 = cyclopropyl) (II). II showed min. inhibitory concentration of ≤0.03, 0.06, 0.013, 0.05, 0.05, 0.20, 0.10, 0.10, 0.013, 0.05, 0.10, 0.10, and 0.39 μg/mL against Escherichia coli NIJH, Shigella flexneri 2A 5503, Proteus vulgaris 08061, Pr. mirabilis IFO-3849, Serratia marcescens 10100, Pseudomonas aeruginosa 32104, Ps. aeruginosa 32121, Xanthomonas maltophilia IID-1275, Staphylococcus aureus 209P, St. epidermidis 56500, Streptococcus pyogenes G-36, Str. faecalis ATCC-19433, and St. aureus 879307, resp. A capsule and a feed dispersant formulation containing II were prepared

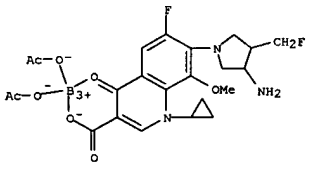
IT **218447-08-6P 218447-10-0P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation of [cis-amino(fluoromethyl)pyrrolidinyl]-1,4-dihydro-4-oxoquinoline-3-carboxylic acid as antibacterial agents)

RN 218447-08-6 CAPLUS
CN Boron, bis(acetato-κO)[1-cyclopropyl-7-[(3S,4S)-3-[[(1,1-dimethylethoxy)carbonyl]amino]-4-(fluoromethyl)-1-pyrrolidinyl]-6-fluoro-1,4-dihydro-8-methoxy-4-(oxo-κO)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)



RN 218447-10-0 CAPLUS
CN Boron, bis(acetato-κO)[7-[(3S,4S)-3-amino-4-(fluoromethyl)-1-pyrrolidinyl]-1-cyclopropyl-6-fluoro-1,4-dihydro-8-methoxy-4-(oxo-κO)-3-quinolinecarboxylato-κO3]-, (T-4)- (9CI) (CA INDEX NAME)

L12 ANSWER 54 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



REFERENCE COUNT: 39 THERE ARE 39 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 55 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1998:684433 CAPLUS
DOCUMENT NUMBER: 129:339856
TITLE: Method, compositions, and aerosol spray containing a polyoxometalate for treating and preventing respiratory viral infections
INVENTOR(S): Schinazi, Raymond F.; Hill, Craig L.
PATENT ASSIGNEE(S): USA
SOURCE: U.S., 18 pp., Cont.-in-part of U.S. Ser. No. 312,561, abandoned.
CODEN: USXXAM
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 2
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 5824706	A	19981020	US 1995-399700	19950303
WO 9609764	A1	19960404	WO 1995-US11961	19950926
W: AM, AT, AU, BB, BG, BR, BY, CA, CH, CN, CZ, DE, DK, EE, ES, FI, GB, GE, HU, IS, JP, KE, KG, KP, KR, KZ, LK, LR, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, TJ, TM				
RW: KE, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
AU 9536366	A1	19960419	AU 1995-36366	19950926
US 6020369	A	20000201	US 1998-111275	19980707
PRIORITY APPLN. INFO.:			US 1994-312561	B2 19940926
			US 1995-399700	A 19950303
			WO 1995-US11961	W 19950926

AB Respiratory viral infections may be effectively prevented or treated by administering an aerosol spray comprising a polyoxometalate to the lungs. (Me3NH)5TaSiW11O40 had a selectivity index greater than 300 when evaluated in HIV-1 acutely infected primary human PBM cells and had no cytotoxicity to uninfected human PBM cells when evaluated up to 100 μM.

IT **131541-70-3**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); THU (Therapeutic use); BIOL (Biological study); USES
(Uses)
(method and aerosol spray containing a polyoxometalate for treating and preventing respiratory viral infections)

RN 131541-70-3 CAPLUS
CN L-Histidine, tetracosa-μ-oxododecaoxo[μ12-[tetrahydroxyborato(5-)-O:O:O:O':O':O':O'':O'':O'':O''':O''':O''']]dodecatungstate(5-) (9CI) (CA INDEX NAME)

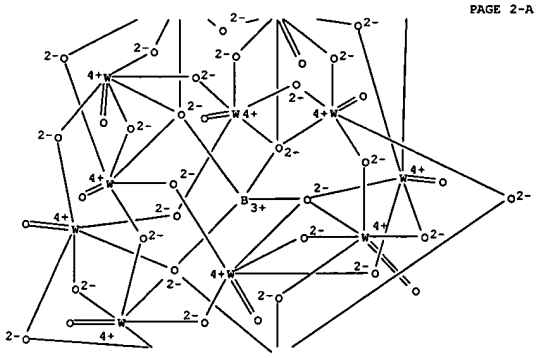
CM 1

CRN 12297-12-0
CMF B O40 W12 . 5 H
CCI CCS

L12 ANSWER 55 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

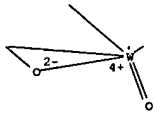
PAGE 1-A

PAGE 2-A



L12 ANSWER 55 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

PAGE 3-A

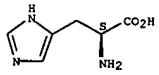


●5 H+

CM 2

CRN 71-00-1
CMF C6 H9 N3 O2

Absolute stereochemistry. Rotation (-).



REFERENCE COUNT: 100 THERE ARE 100 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 56 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1997:454680 CAPLUS
DOCUMENT NUMBER: 127:183157
TITLE: Chemical studies on the nonlinear optics of coordination compounds
AUTHOR(S): You, Xiao-Zeng
CORPORATE SOURCE: Coordination Chemistry Institute, State Key Laboratory
Coordination Chemistry, Center Advanced Studies Sciences and Technology Mirostructures, Nanjing University, Nanjing, 210093, Peop. Rep. China
SOURCE: Journal of Photochemistry and Photobiology, A: Chemistry (1997), 106(1-3), 85-90
CODEN: JPPCEJ; ISSN: 1010-6030
PUBLISHER: Elsevier
DOCUMENT TYPE: Journal
LANGUAGE: English

AB There has been much progress in chemical studies on nonlinear optics organic, polymer, inorg. and organometallic solid materials, but less for coordination compds. This paper present some of our recent research with emphasis on the possibility of incorporation advantages of both organic (high nonlinear optical effect) and inorg. (stable large crystal) parts into the form of coordination compds. (include organic-inorg. salts) for nonlinear optical applications.

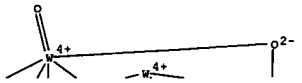
IT **194150-79-3 194150-80-6**
RL: PRP (Properties)
(nonlinear optical properties of coordination compds.)

RN 194150-79-3 CAPLUS
CN Tungstate(5-), tetracosa-μ-oxododecaoxo[μ12-[tetrahydroxyborato(5-)-κO:κO:κO:κO':κO':κO':κO''::.kappa.O'':κO'':κO''':κO''':κO''']]dodeca-, pentahydrogen, compd. with 1-methyl-2-pyrrolidinone (1:5) (9CI) (CA INDEX NAME)

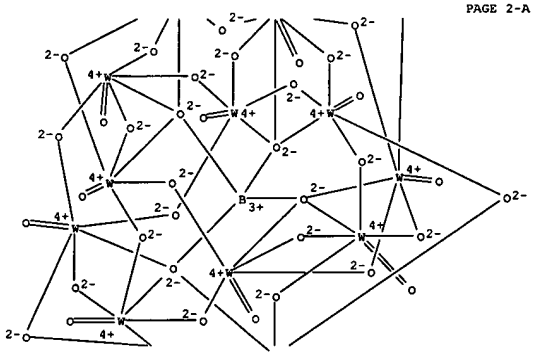
CM 1

CRN 12297-12-0
CMF B O40 W12 . 5 H
CCI CCS

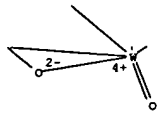
PAGE 1-A



PAGE 2-A



PAGE 3-A



●5 H^+

CM 2

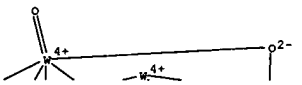
CRN 872-50-4
CMF C5 H9 N O

RN 194150-80-6 CAPLUS
CN Tungstate(5-), tetracosa-μ-oxododecaoxo[μ12-[tetrahydroxyborato(5-)-κO:κO:κO:κO':κO':κO'':κO'':κO''':κO''':κO''']]dodeca-, pentahydrogen, compd. with 1-methyl-2-pyrrolidinone (1:5), dihydrate (9CI)
(CA INDEX NAME)

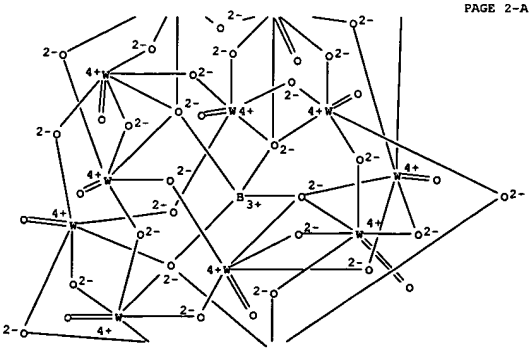
CM 1

CRN 12297-12-0
CMF B O40 W12 . 5 H
CCI CCS

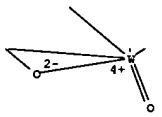
PAGE 1-A



PAGE 2-A



PAGE 3-A



●5 H^+

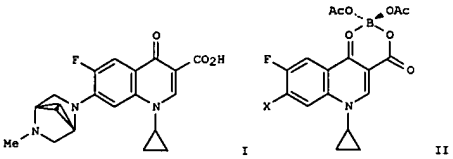
CM 2

CRN 872-50-4
CMF C5 H9 N O

L12 ANSWER 57 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1997:388730 CAPLUS
DOCUMENT NUMBER: 127:5021
TITLE: Process for preparation of 1-cyclopropyl-6-fluoro-1,4-dihydro-7-[(1S,4S)-5-methyl-2,5-diazabicyclo[2.2.1]hept-2-yl]-4-oxo-3-quinolinecarboxylic acid [danofloxacin] and its salts
INVENTOR(S): Picornell Darder, Carlos; Gonzalez Harnandez, Pedro; Salas Gonzalez, M. Luisa
PATENT ASSIGNEE(S): Quimica Sintetica, S.A., Spain
SOURCE: Span., 7 pp.
CODEN: SPXXAD
DOCUMENT TYPE: Patent
LANGUAGE: Spanish
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

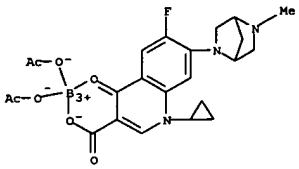
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
ES 2092963	A1	19961201	ES 1995-782	19950412
ES 2092963	B1	19971216		
FI 9601583	A	19961013	FI 1996-1583	19960411
NO 9601442	A	19961014	NO 1996-1442	19960411
AT 9600653	A	19980915	AT 1996-653	19960411
PRIORITY APPLN. INFO.:			ES 1995-782 A	19950412

OTHER SOURCE(S): CASREACT 127:5021
GI



AB Title compound I and salts are prepared in 3 steps. Treatment of a C1-4 alkyl 1-cyclopropyl-6-fluoro-7-chloro-1,4-dihydro-4-oxo-3-quinolinecarboxylate with boric acid in the presence of Ac2O and catalytic Zn gives the boron chelate intermediate II [X = Cl]. Treatment of the latter with (1S,4S)-5-methyl-2,5-diazabicyclo[2.2.1]heptane or its addition salts, optionally in the presence of a base, gives the intermediate II [X = 7-[(1S,4S)-5-methyl-2,5-diazabicyclo[2.2.1]hept-2-yl]]. Finally, hydrolysis of this chelate (e.g., with aqueous NaOH) gives I. In 2 examples, a combination of the 2nd and 3rd steps gave I in 92-94% yield.
IT **190132-00-4P**
RL: IMF (Industrial manufacture); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(intermediate; preparation of diazabicycloheptyloxoquinolinecarboxylic acid

L12 ANSWER 57 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
deriv. via boron chelates)
RN 190132-00-4 CAPLUS
CN Boron, bis(acetato-κO)[1-cyclopropyl-6-fluoro-1,4-dihydro-7-(5-methyl-2,5-diazabicyclo[2.2.1]hept-2-yl)-4-(oxo-κO)-3-quinolinecarboxylato-κO3]-, [T-4-(1S)]- (9CI) (CA INDEX NAME)



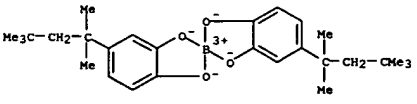
L12 ANSWER 58 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1997:334873 CAPLUS
DOCUMENT NUMBER: 126:310442
TITLE: Electrophotographic toner containing complex salts
INVENTOR(S): Nagatsuka, Takayuki; Tanaka, Katsuhiko
PATENT ASSIGNEE(S): Canon Kk, Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 17 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 09068824	A2	19970311	JP 1995-245102	19950831
JP 3323710	B2	20020909		
PRIORITY APPLN. INFO.:			JP 1995-245102	19950831

AB The toner contains fine particles of resin, inorg. oxide fine powders, and complex salts which satisfy the relations Vc ≥ 0.3 nm2 and Va > Vc (Va and Vc = volume of anion and cation, resp.) and have 5- or 6-membered ring as the metal coordination site. The fine particles of resin may contain magnetic powder which shows variation coefficient of the particle size ≤35% and satisfies the relation: (-7/3) + r + 45 ≤ MT ≤ (-7/3) + r + 75 [MT = content of the magnetic powder; r = weight average particle size of the toner (μm)] for uniform dispersion. The toner shows good environmental stability and provide images excellent in uniformity at the highlighted area.
IT **189263-24-9**
RL: PRP (Properties); TEM (Technical or engineered material use); USES (Uses)
(electrophotog. toner containing complex salts composed of volume-defined anion and cation to improve environmental stability)
RN 189263-24-9 CAPLUS
CN 1H-Imidazolium, 2-methyl-1,3-bis(phenylmethyl)-, (T-4)-bis[4-(1,1,3,3-tetramethylbutyl)-1,2-benzenediolato(2-)-κO,κO']borate(1-) (9CI) (CA INDEX NAME)

CM 1

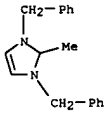
CRN 189263-22-7
CMF C28 H40 B O4
CCI CCS



CM 2

L12 ANSWER 58 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CRN 101554-35-2
CMF C18 H19 N2



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

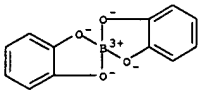
L12 ANSWER 59 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1996:381137 CAPLUS
DOCUMENT NUMBER: 125:128193
TITLE: Complexes of boron with catechol: the x-ray crystal structure of Meulenhoff's salt, 2[(pyH)[B(cat)2]·H2cat]·H2O
AUTHOR(S): Griffith, William P.; White, Andrew J. P.; Williams, David J.
CORPORATE SOURCE: Inorganic and Chemical Crystallographic Research Laboratories, Imperial Coll. of Sci., Technol. and Med., London, SW7 2AY, UK
SOURCE: Polyhedron (1996), 15(17), 2835-2839
CODEN: PLYHDE; ISSN: 0277-5387
PUBLISHER: Elsevier
DOCUMENT TYPE: Journal
LANGUAGE: English

AB Meulenhoff's salt, 2[(pyH)[B(cat)2]·H2cat]·H2O, was prepared by the original literature method and its x-ray crystal structure determined
Crystals are monoclinic, space group P21/n, with a 24.12(2), b 7.090(8), c
25.04(2) Å, and β 96.52(7)°; Z = 4, dc = 1.33; R = 0.066,
Rw = 0.055. The stoichiometry is confirmed as 1:1:3 pyridine:B:catechol, though only two of the catechol units are bound directly to the B center. The IR and Raman vibrational spectra of the solid and the 1H, 13C{1H} and 11B{1H} NMR spectra of the complex in solution are reported, and suggest that
the solid state structure of the anion is maintained in solution

IT **179122-55-5**
RL: PRP (Properties)
(crystal structure of)
RN 179122-55-5 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 1,2-benzenediol and pyridine, hydrate (2:2:2:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



CM 2

CRN 120-80-9
CMF C6 H6 O2

L12 ANSWER 59 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

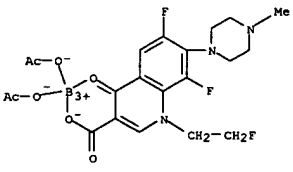
CM 3

CRN 110-86-1
CMF C5 H5 N

L12 ANSWER 60 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1996:269580 CAPLUS
DOCUMENT NUMBER: 125:10649
TITLE: Studies on synthesis of a fluoroquinolone antimicrobial agent fleroxacin
AUTHOR(S): Wang, Erhua; Jin, Rong
CORPORATE SOURCE: Medical and Chem. Inst., China Pharmaceutical Univ., Nanjing, 210009, Peop. Rep. China
SOURCE: Zhongguo Yaoke Daxue Xuebao (1995), 26(6), 321-3
CODEN: ZHYXE9; ISSN: 1000-5048
PUBLISHER: Zhongguo Yaoke Daxue
DOCUMENT TYPE: Journal
LANGUAGE: Chinese

AB An improved synthesis of antibacterial agent fleroxacin, starting from 2,3,4-trifluoronitrobenzene via reduction, condensation-cyclization, fluoroethylation, chelation, piperazination, and hydrolysis was described.
The reactions of all steps were carried out under moderate condition with the overall yield of 40%.

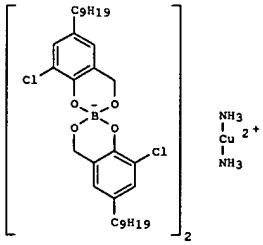
IT **176982-64-2P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(synthesis of fluoroquinolone antimicrobial agent fleroxacin from trifluoronitrobenzene)
RN 176982-64-2 CAPLUS
CN Boron, bis(acetato-O)[6,8-difluoro-1-(2-fluoroethyl)-1,4-dihydro-7-(4-methyl-1-piperazinyl)-4-oxo-3-quinolincarboxylato-O3,O4]-, (T-4)- (9CI) (CA INDEX NAME)



L12 ANSWER 61 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1996:163906 CAPLUS
DOCUMENT NUMBER: 124:202593
TITLE: Novel co-ordinated metal/boron compounds as biocides, their methods of synthesis, their use and their formulation
INVENTOR(S): Maynard, Nigel Paul
PATENT ASSIGNEE(S): N. Z.
SOURCE: PCT Int. Appl., 38 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9531425	A1	19951123	WO 1995-NZ38	19950508
W: AM, AT, AU, BB, BG, BR, BY, CA, CH, CN, CZ, DE, DK, EE, ES, FI, GB, GE, HU, IS, JP, KE, KG, KP, KR, KZ, LK, LR, LT, LU, LV, MD, MG, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, TJ, TM, TT				
RW: KE, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
AU 9525393	A1	19951205	AU 1995-25393	19950508
ZA 9503931	A	19960117	ZA 1995-3931	19950515
PRIORITY APPLN. INFO.:			NZ 1994-260530 A	19940516
			WO 1995-NZ38 W	19950508

OTHER SOURCE(S): MARPAT 124:202593
GI



I

AB The invention includes the synthesis of multivalent metal complexes with B
containing complex organic anions, e.g., I. The complexes include or may include

L12 ANSWER 61 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

ligands. Complexes capable of being so synthesized are also disclosed as is their use as biocides which fix owing to low aq. soly. Intermediates of the process or complexes of the process (after heating or when provided with ligands low on the series) are also useful to strip multivalent metal cations or ligand forming compds. from an aq. phase (usually by the movement of the complex with the attached multivalent metal cations or ligands into an org. solvent phase), thereby completing the synthesis.

IT **174459-03-1P**

RL: AGR (Agricultural use); BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation as biocides)

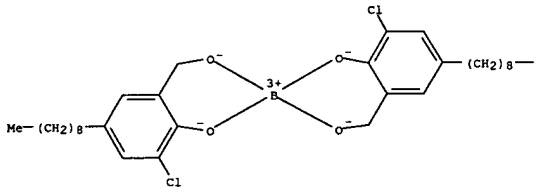
RN 174459-03-1 CAPLUS

CN Copper(2+), bis(pyridine)-, bis[(T-4)-bis[3-chloro-2-hydroxy-5-nonylbenzenemethanolato(2-)-O,O']borate(1-)] (9CI) (CA INDEX NAME)

CM 1

CRN 174459-00-8
CMF C32 H46 B Cl2 O4
CCI CCS

PAGE 1-A



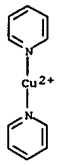
PAGE 1-B

—Me

CM 2

CRN 23236-29-5
CMF C10 H10 Cu N2
CCI CCS

L12 ANSWER 61 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



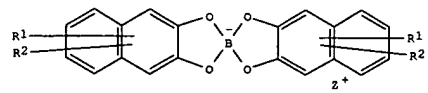
L12 ANSWER 62 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1996:67579 CAPLUS
DOCUMENT NUMBER: 124:215974
TITLE: Electrostatographic developer toner with good charging properties
INVENTOR(S): Tanaka, Katsuhiko; Nagatsuka, Takayuki; Ichikawa, Yasuhiro; Takahashi, Toshihiko
PATENT ASSIGNEE(S): Canon Kk, Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 18 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

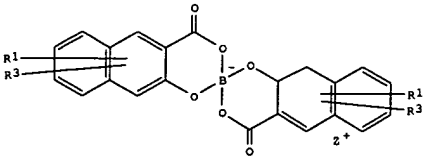
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 07295296	A2	19951110	JP 1994-110243	19940427
JP 3074589	B2	20000807		
PRIORITY APPLN. INFO.:			JP 1994-110243	19940427

OTHER SOURCE(S): MARPAT 124:215974

GI



I



II

AB The toner contains ≥1 I and/or II (R1-2 = H, halo, alkyl, alkoxy, acyl, ester; R3 = halo, alkyl, alkoxy, acyl, ester; R1-3 may be substituted with aryl group). The toner may contain an inorg. oxide with electronegativity of metal ion 10-15. The toner showed good charging properties and repeating durability.

IT **174188-61-5**

RL: TEM (Technical or engineered material use); USES (Uses)
(electrostatog. developer toner containing boron compound charge-controlling agent)

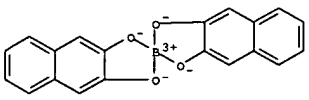
RN 174188-61-5 CAPLUS

CN Borate(1-), bis[2,3-naphthalenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

L12 ANSWER 62 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CM 1

CRN 47422-29-7
CMF C20 H12 B O4 . H
CCI CCS



CM 2

CRN 110-86-1
CMF C5 H5 N

L12 ANSWER 63 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

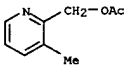
ACCESSION NUMBER: 1995:742629 CAPLUS
DOCUMENT NUMBER: 123:143645
TITLE: Improvements in the object of patent 9,102,594 for a process for the preparation of pyridine derivatives.
INVENTOR(S): Palomo Coll, Alberto
PATENT ASSIGNEE(S): Centro Genesis para la Investigacion S.L., Spain
SOURCE: Span., 15 pp.
CODEN: SPXXAD
DOCUMENT TYPE: Patent
LANGUAGE: Spanish
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
ES 2058027	A1	19941016	ES 1993-240	19930209
ES 2058027	B1	19950516		
PRIORITY APPLN. INFO.:			ES 1991-2590 A	19911121

OTHER SOURCE(S): CASREACT 123:143645; MARPAT 123:143645
GI For diagram(s), see printed CA Issue.
AB In the title-referenced process for the preparation of pyridine N-oxides I and pyridinemethanols II [R2, R3 = H, Me, OMe; R4 = CH2CF3, Et, iso-Pr, (CH2)3OMe], the improvement is characterized by reaction, in a solvent, of the reactive intermediate salts III or IV [D = variable-valence element; a, b = 0 or integer, with (a+b)>0] with a corresponding alc. derivative D(OR4)c-M+ [V; M = Lewis acid, protonated organic base, Si derivative]. Elements functioning as D include Pd, P, S, B, and Br. The method is applicable to preparation of intermediates for antiulcer agents, especially lansoprazole. For example, 0.4 mL PBr3 was heated with 5 mL CF3CH2OH at 70° until formed HBr was eliminated (solution A). Meanwhile, 3 g 85% KOH was added to 10 mL CF3CH2OH at 5°, with the temperature rising to 20-25° (solution B). Upon dissoln. of KOH, solution A was added to solution B, followed by 2.5 g 2,3-dimethyl-4-nitropyridine N-oxide. Refluxing of the mixture for 2 h 25 min gave complete reaction to a single product by TLC, namely 2,3-dimethyl-4-(2,2,2-trifluoroethoxy)pyridine N-oxide (VI). Addnl. examples show preparation of VI using SCl2 or Br2 in place of PBr3. Evidence for the existence of the denitro-supercations in III and IV, and the superanions in V, is described.
IT **166521-88-6P 166521-89-7P**
RL: PNU (Preparation, unclassified); PREP (Preparation)
(reactive intermediate; improved preparation of alkoxypyridine derivs. via alkoxydenitration of nitropyridines with supersalts)
RN 166521-88-6 CAPLUS
CN Borate(1-), (nitrito-O)tris(2,2,2-trifluoroethanolato)-, (T-4)-, salt with 2,3-dimethylpyridine 1-oxide (1:1) (9CI) (CA INDEX NAME)

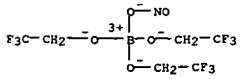
CM 1

L12 ANSWER 63 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



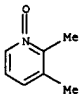
L12 ANSWER 63 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CRN 166521-87-5
CMF C6 H6 B F9 N O5
CCI CCS



CM 2

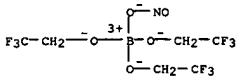
CRN 166521-76-2
CMF C7 H9 N O
CCI RIS



RN 166521-89-7 CAPLUS
CN Borate(1-), (nitrito-O)tris(2,2,2-trifluoroethanolato)-, (T-4)-, salt with (3-methyl-2-pyridinyl)methyl acetate (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 166521-87-5
CMF C6 H6 B F9 N O5
CCI CCS



CM 2

CRN 166521-79-5
CMF C9 H11 N O2
CCI RIS

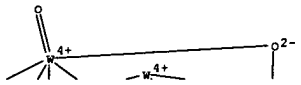
L12 ANSWER 64 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1995:206843 CAPLUS
DOCUMENT NUMBER: 122:281395
TITLE: Relationship of the molecular size and charge density of polyoxometalates to their anti-gp120-CD4-binding activity
AUTHOR(S): Judd, D. A.; Schinazi, R. F.; Hill, C. L.
CORPORATE SOURCE: Dep. Chem., Emory Univ., Atlanta, GA, 30322, USA
SOURCE: Antiviral Chemistry & Chemotherapy (1994), 5(6), 410-14
CODEN: ACCHEH; ISSN: 0956-3202
PUBLISHER: Blackwell
DOCUMENT TYPE: Journal
LANGUAGE: English
AB The ability of several classes of polyoxometalates to inhibit the interaction between HIV-1 gp120 and CD4 was assessed. No clear relation was found between binding inhibition and the neg. charge d. on the anion portion of the polyoxometalate. However, a weak correlation was found with mol. size. There was a mol. weight threshold of 3800 g mol-1 above which no significant increase in potency was gained; the binding inhibition was nearly quant. above this mol. weight
IT **131541-70-3**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); PRP (Properties); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(relationship of mol. size and charge d. of polyoxometalates to their anti-gp120-CD4-binding activity)
RN 131541-70-3 CAPLUS
CN L-Histidine, tetracosa-μ-oxododecaoxo[μ12-[tetrahydroxyborato(5-)-O:O:O:O':O':O':O'':O'':O'':O''':O''':O''']]dodecatungstate(5-) (9CI) (CA INDEX NAME)

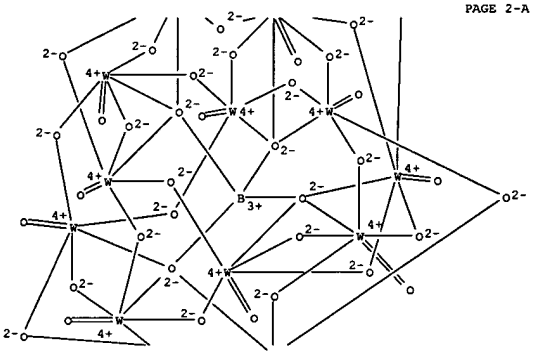
CM 1

CRN 12297-12-0
CMF B O40 W12 . 5 H
CCI CCS

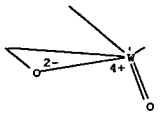
PAGE 1-A



PAGE 2-A



PAGE 3-A

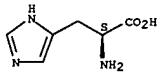


●5 H^+

CM 2

CRN 71-00-1
CMF C6 H9 N3 O2

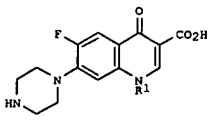
Absolute stereochemistry. Rotation (-).



ACCESSION NUMBER: 1995:182998 CAPLUS
DOCUMENT NUMBER: 122:81399
TITLE: Process for preparing 1-substituted-6-fluoro-4-oxo-7-(1-piperazinyl)-1,4-dihydroquinoline-3-carboxylic acid, a novel intermediate useful in said process, and
a process for preparing said intermediate
INVENTOR(S): Zupancic, Natasa; Barbo, Martin; Sket, Boris; Zupet, Pavel
PATENT ASSIGNEE(S): Slovenia
SOURCE: Can. Pat. Appl., 14 pp.
CODEN: CPXXEB
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
CA 2111181	AA	19940612	CA 1993-2111181	19931210
HU 75319	A2	19970528	HU 1993-2940	19931018
PL 173784	B1	19980430	PL 1993-301045	19931112
CZ 284715	B6	19990217	CZ 1993-2643	19931206
LT 3084	B	19941125	LT 1993-1558	19931207
AT 9302497	A	19960315	AT 1993-2497	19931210
AT 401648	B	19961025		
LV 10863	B	19960820	LV 1993-1317	19931210
RU 2127270	C1	19990310	RU 1993-54527	19931210
PRIORITY APPLN. INFO.:			SI 1992-377 A	19921211

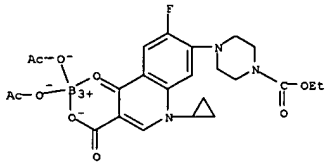
OTHER SOURCE(S): CASREACT 122:81399; MARPAT 122:81399
GI



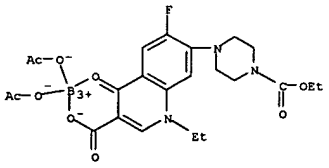
I

AB A process for preparing 1-substituted-6-fluoro-4-oxo-7-(1-piperazinyl)-1,4-dihydroquinoline-3-carboxylic acids I (R2 = alkyl, cycloalkyl or 2,4-difluorophenyl) was disclosed; I are prepared by hydrolysis of novel boron diacetate precursors. II are useful for the treatment of inflammatory diseases. It is obtained by a nucleophilic substitution of the halo atom in 7-position in the compound of the formula III.
IT **158964-91-1P 158964-92-2P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation of (fluoro)oxo(piperazinyl)quinolinecarboxylates from boron
diacetate intermediates)
RN 158964-91-1 CAPLUS

CN Boron, bis(acetato-O)[1-cyclopropyl-7-[4-(ethoxycarbonyl)-1-piperazinyl]-6-fluoro-1,4-dihydro-4-oxo-3-quinolinecarboxylato-O3,O4]-, (T-4)- (9CI) (CA
INDEX NAME)



RN 158964-92-2 CAPLUS
CN Boron, bis(acetato-O)[7-[4-(ethoxycarbonyl)-1-piperazinyl]-1-ethyl-6-fluoro-1,4-dihydro-4-oxo-3-quinolinecarboxylato-O3,O4]-, (T-4)- (9CI) (CA
INDEX NAME)

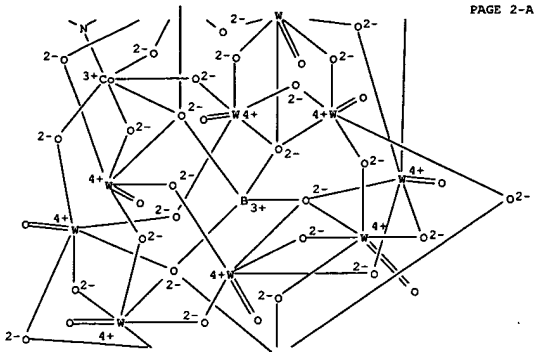


L12 ANSWER 66 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1994:547566 CAPLUS
DOCUMENT NUMBER: 121:147566
TITLE: NMR spectra of 4,4'-bipyridyl, pyrazine, and ethylenediamine coordinated to undecatungstocobalto(III)silicate and -borate anions. Identification of 1:1 and dumbbell-shaped 1:2 complexes
AUTHOR(S): Park, Jeongmin; Ko, Moonhee; So, Hyunsoo
CORPORATE SOURCE: Dep. Chem., Sogang Univ., Seoul, 121-742, S. Korea
SOURCE: Bulletin of the Korean Chemical Society (1993), 14(6), 759-62
CODEN: BKCSDE; ISSN: 0253-2964
DOCUMENT TYPE: Journal
LANGUAGE: English
AB The reaction of Cs2[SiW11O39Co(H2O)] (I) and K6[BW11O39Co(H2O)] (II) with en, 4,4'-bipyridine and pyrazine was studied by NMR. 4,4'-Bpy forms 1:1 and 2:1 complexes with I and a 1:1 complex with II. En reduces Co(III) in I but forms 1:1 and 2:1 complexes with II. Pyrazine forms 1:1 and 2:1 complexes with II and 1 1:1 complex with I.
IT **157177-67-8P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 157177-67-8 CAPLUS
CN Tungstate(6-), tetracosa-μ-oxoundecaoxo[(pyrazine-N1)cobaltate][μ12-[tetrahydroxyborato(5-)-O:O:O:O':O':O':O'':O'':O'':O''':O''':O''']]undeca- (9CI) (CA INDEX NAME)

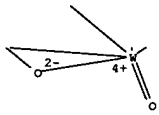
* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

L12 ANSWER 66 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

PAGE 2-A



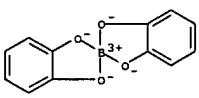
PAGE 3-A



L12 ANSWER 67 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1994:485097 CAPLUS
DOCUMENT NUMBER: 121:85097
TITLE: Microfiltration membranes with surface charge
AUTHOR(S): Grachek, V. I.; Chekhovich, L. P.; Antomonov, V. A.
CORPORATE SOURCE: Inst. Fiz. Org., Belarus
SOURCE: Vestsi Akademii Navuk Belarusi, Seryya Khimichnykh Navuk (1993), (4), 70-3
CODEN: VAKNEK; ISSN: 1025-5567
DOCUMENT TYPE: Journal
LANGUAGE: Russian
AB Improved polyamide-6 membranes with pos. zeta potential with pH from 3.7 to 10.5 have been prepared by treatment with compds. containing cationic functional groups and crosslinking agents (epoxides). The porous structure and performance characteristics of modified membranes were not affected by modification.
IT **144672-73-1**
RL: USES (Uses)
(polyamide membranes modified by, electrokinetic potential and performance characteristics of)
RN 144672-73-1 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H+

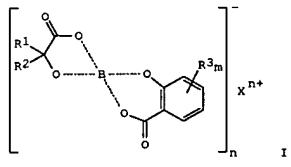
CM 2

CRN 504-29-0
CMF C5 H6 N2

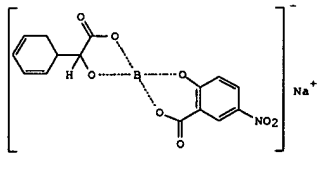
L12 ANSWER 68 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1994:257410 CAPLUS
DOCUMENT NUMBER: 120:257410
TITLE: Electrophotographic toners using boron complex charge-controlling agent
INVENTOR(S): Hasegawa, Junko; Takeda, Teruichi; Ebisawa, Makoto
PATENT ASSIGNEE(S): Japan Carlit Co Ltd, Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 9 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 05265257	A2	19931015	JP 1992-91563	19920318
PRIORITY APPLN. INFO.:			JP 1992-91563	19920318

OTHER SOURCE(S): MARPAT 120:257410
GI



I



II

AB The title toners contain I [R1, R2 = H, alkyl, (substituted) aromatic ring (including condensed ring); R3 = H, alkyl, alkoxy, halo, NO2, sulfone, amide, sulfamide, aromatic ring (including condensed ring); Xn+ = cation; m = 1-4] as a charge-controlling agent in a polyester resin. The complex is colorless, shows good compatibility with polyester resins, and the toners exhibit improved environmental stability and provide high color quality images. Thus, a polyester resin, carbon black, and II were kneaded and pulverized to give a toner, which was mixed with an Fe powder to give a developer.

L12 ANSWER 68 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

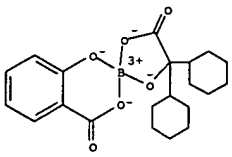
IT **154507-78-5 154507-80-9**
RL: USES (Uses)
(charge-controlling agent, electrophotog. toner containing)

RN 154507-78-5 CAPLUS

CN Borate(1-), [α-cyclohexyl-α-hydroxycyclohexaneacetato(2-)][2-hydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 1-methyl-1H-pyrrole (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 154507-77-4
CMF C21 H26 B O6 . H
CCI CCS



CM 2

CRN 96-54-8
CMF C5 H7 N

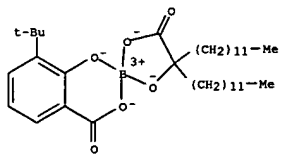
RN 154507-80-9 CAPLUS

CN Piperidinium, 1,1-dimethyl-, (T-4)-[3-(1,1-dimethylethyl)-2-hydroxybenzoato(2-)-O1,O2][2-dodecyl-2-hydroxytetradecanoato(2-)-O1,O2]borate(1-) (9CI) (CA INDEX NAME)

CM 1

CRN 154507-79-6
CMF C37 H62 B O6
CCI CCS

L12 ANSWER 68 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



CM 2

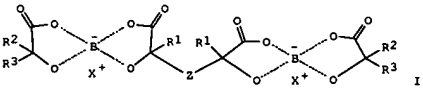
CRN 15302-91-7
CMF C7 H16 N

L12 ANSWER 69 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

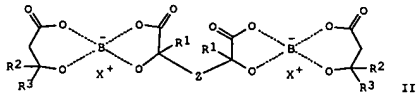
ACCESSION NUMBER: 1994:178138 CAPLUS
DOCUMENT NUMBER: 120:178138
TITLE: Electrophotographic toners using boron complex as charge-controlling agent
INVENTOR(S): Hasegawa, Junko; Takeda, Teruichi; Ebisawa, Makoto
PATENT ASSIGNEE(S): Japan Carlit Co Ltd, Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 10 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 05165256	A2	19930702	JP 1991-351823	19911216
PRIORITY APPLN. INFO.:			JP 1991-351823	19911216

GI



I



II

AB The title toners contain, as a charge-controlling agent, a B complex I or II [R1-3 = H, alkyl, alkenyl, (substituted) aromatic ring or condensed ring; Z = (substituted) aromatic ring or condensed ring; X+ = cation]. The charge-controlling agent is colorless and the toners show sharp and uniform triboelec. charge distribution and good environmental stability and provide high-quality color images. Thus, a polystyrene resin, carbon black, and I (R1 = R2 = H; R3 = Ph; Z = p-phenylene; X+ = Li+) were kneaded and pulverized to give a toner, which was mixed with an Fe powder to give a developer.

IT **152576-36-8 152576-39-1**
RL: USES (Uses)
(charge-controlling agent, electrophotog. photoreceptor containing)

RN 152576-36-8 CAPLUS

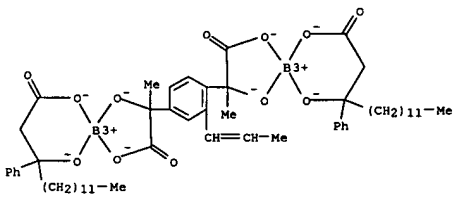
CN Piperidinium, 1,1-dimethyl-, [μ-[α,α'-dihydroxy-α,α'-dimethyl-3-(1-propenyl)-1,4-benzenediacetato(4-)]]bis[β-dodecyl-β-hydroxybenzenepropanoato(2-)-Oα,Oβ]diborate(2-) (2:1) (9CI) (CA INDEX NAME)

CM 1

CRN 152576-35-7

L12 ANSWER 69 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CMF C57 H78 B2 O12
CCI CCS



CM 2

CRN 15302-91-7
CMF C7 H16 N

RN 152576-39-1 CAPLUS

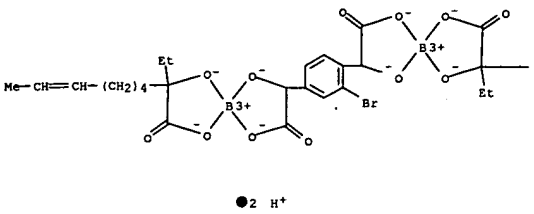
CN Borate(2-), [μ-[2-bromo-α,α'-dihydroxy-1,4-benzenediacetato(4-)]]bis[2-ethyl-2-hydroxy-7-nonenoato(2-)-O1,O2]di-, dihydrogen, compd. with 1-methyl-1H-pyrrole (1:2) (9CI) (CA INDEX NAME)

CM 1

CRN 152576-38-0
CMF C32 H41 B2 Br O12 . 2 H
CCI CCS

L12 ANSWER 69 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

PAGE 1-A



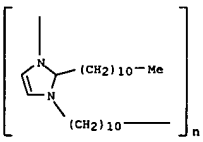
●2 H^+

PAGE 1-B

CM 2

CRN 96-54-8
CMF C5 H7 N

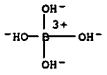
L12 ANSWER 70 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

CM 2

CRN 15390-83-7
CMF B H4 O4
CCI CCS

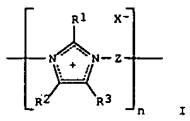


L12 ANSWER 70 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1994:90789 CAPLUS
DOCUMENT NUMBER: 120:90789
TITLE: Charge-controlling agents and electrophotographic toners using them
INVENTOR(S): Nakanishi, Hideo; Ikeda, Hiroyuki
PATENT ASSIGNEE(S): Sanyo Chemical Ind Ltd, Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 9 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 05165258	A2	19930702	JP 1991-352241	19911213
JP 08012486	B4	19960207		
PRIORITY APPLN. INFO.:			JP 1991-352241	19911213

GI



AB The charge-controlling agents are polymer quaternary salts I (R1 = H, C1-17 hydrocarbon; R2, R3 = H, C1-8 hydrocarbon, R2 and R3 may form an aromatic ring; Z = C1-12 alkylene group which may have ether bond; X- = anion; n = 2-100), and the toners comprise a binder resin, a colorant, and the agent. The toners show good thermal resistance and stable chargeability under varied environmental conditions and prevent generation of bad smell on heating. Thus, Himer GRX2500 (styrene-acrylic copolymer), Regal 330R (carbon black), I [R1-3 = H; Z = (CH2)6; X- = 1-naphthalenesulfonate ion], and Viscol 550P (polypropylene wax) were kneaded and pulverized to give a toner, which was mixed with a ferrite carrier to give a developer.
IT **152584-99-1**
RL: USES (Uses)
(charge-controlling agent, electrophotog. developer toner containing)
RN 152584-99-1 CAPLUS
CN Poly[(2-undecyl-1H-imidazolium-1,4-diyl)-1,10-decanediyl tetrahydroxyborate(1-)] (9CI) (CA INDEX NAME)

CM 1

CRN 152584-98-0
CMF (C24 H45 N2)n
CCI PMS

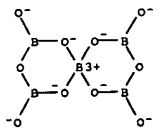
L12 ANSWER 71 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1993:661384 CAPLUS
DOCUMENT NUMBER: 119:261384
TITLE: Clathrates with three-dimensional host structures of hydrogen-bonded pentaborate [B5O6(OH)4]- ions: pentaborates with the cations NMe4+, NEt4+, NPhMe3+ and pipH+ (pipH+ = piperidinium)
AUTHOR(S): Wiebcke, Michael; Freyhardt, Clemens C.; Felsche, Juergen; Engelhardt, Guenter
CORPORATE SOURCE: Fak. Chem., Univ. Konstanz, Konstanz, D-78434, Germany
SOURCE: Zeitschrift fuer Naturforschung, B: Chemical Sciences (1993), 48(7), 978-85
CODEN: ZNBSEN; ISSN: 0932-0776
DOCUMENT TYPE: Journal
LANGUAGE: English
AB X-ray structure analyses of crystalline [NMe4][B5O6(OH)4].nH2O (n = 0.25-0.50) (1), [NEt4][B5O6(OH)4] (2), [NPhMe3][B5O6(OH)4] (3), and [pipH][B5O6(OH)4] (4) reveal that these materials are novel clathrates with closely related 3-dimensional host structures built up of H-bonded oligomeric [B5O6(OH)4]-. The organic cations and H2O mols. (in 1) occupy as guest species large straight channel-like voids of nearly rectangular cross section. 1 Crystallizes as monoclin., space group P21/c, Z = 4; 2, 3 and 4, which possess the same host-structure topol., crystallize as triclinic, space group P.hivin.1, Z = 2. 11B MAS NMR spectra allow the detection of small angular distortions in the [B5O6(OH)4]- caused by the specific H bonding in the host frameworks. Upon heating the compds. on a thermobalance in a dynamic inert gas atmospheric dehydration occurs at temps. of 563 K (1), 543 K (2), 558 K (3) and 523 K (4) before degradation of the organic cations starts at temps. of 633 K (1), 623 K (2), 623 K (3) and 613 K (4).
IT **12548-84-4P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation and crystal structure and boron-11 NMR spectrum and thermal dehydration and decomposition of)
RN 12548-84-4 CAPLUS
CN Borate(5-), bis[μ-oxotetraoxodiborato(4-)]-, (T-4)-, pentahydrogen, compd. with piperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 92258-67-8
CMF B5 O10 . 5 H
CCI CCS

L12 ANSWER 71 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



●5 H+

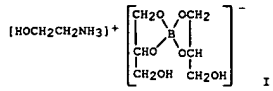
CM 2

CRN 110-89-4
CMF C5 H11 N

L12 ANSWER 72 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1993:429575 CAPLUS
DOCUMENT NUMBER: 119:29575
TITLE: Antistatic thermoplastic resin laminates
INVENTOR(S): Kuze, Katsuro; Tahoda, Tadashi
PATENT ASSIGNEE(S): Toyobo Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 8 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 04311735	A2	19921104	JP 1991-77792	19910410
PRIORITY APPLN. INFO.:			JP 1991-77792	19910410

OTHER SOURCE(S): MARPAT 119:29575
GI

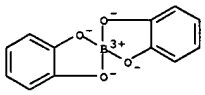


I

AB The title laminates comprise antistatic agent layers of charge-transfer complexes comprising nonionic amines and semipolar borates (e.g., I) on ≥1 side of a thermoplastic resin film (e.g., PET).
IT **147390-96-3 147390-98-5**
RL: USES (Uses)
(antistatic agents, for thermoplastic films)
RN 147390-96-3 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methylmorpholine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H+

L12 ANSWER 72 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

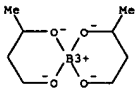
CM 2

CRN 109-02-4
CMF C5 H11 N O

RN 147390-98-5 CAPLUS
CN Borate(1-), bis[1,3-butanediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 147390-97-4
CMF C8 H16 B O4 . H
CCI CCS



● H+

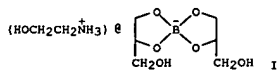
CM 2

CRN 110-86-1
CMF C5 H5 N

L12 ANSWER 73 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1993:214286 CAPLUS
DOCUMENT NUMBER: 118:214286
TITLE: Antistatic agents for plastics, fibers, or paper
INVENTOR(S): Kuze, Katsuro; Tahoda, Tadashi; Hamanaka, Hiroyoshi
PATENT ASSIGNEE(S): Toyobo Co., Ltd., Japan; Boron International K.K.
SOURCE: Jpn. Kokai Tokkyo Koho, 10 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 04288390	A2	19921013	JP 1991-15536	19910206
JP 2968061	B2	19991025		
PRIORITY APPLN. INFO.:			JP 1991-15536	19910206

GI

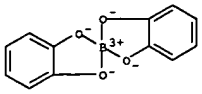


I

AB The title agents comprise the reaction products of nonionic compds. containing
one basic N and B compds. with a specified structure. Thus, a polyester film was coated with an aqueous solution containing I to give a 0.5-μm coat with
surface resistivity 3.8 + 1010 Ω, vs. >1015 for the noncoated film.
IT **147390-96-3 147390-98-5**
RL: USES (Uses)
(antistatic agents, for plastic films)
RN 147390-96-3 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methylmorpholine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H^+

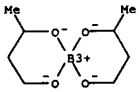
CM 2

CRN 109-02-4
CMF C5 H11 N O

RN 147390-98-5 CAPLUS
CN Borate(1-), bis[1,3-butanediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 147390-97-4
CMF C8 H16 B O4 . H
CCI CCS



● H^+

CM 2

CRN 110-86-1
CMF C5 H5 N

ACCESSION NUMBER: 1992:652822 CAPLUS
DOCUMENT NUMBER: 117:252822
TITLE: Borate-containing polyamide composition for membrane filters
INVENTOR(S): Grachek, V. I.; Artamonov, V. A.; Chekhovich, L. P.
PATENT ASSIGNEE(S): Institute of Physical-Organic Chemistry, Academy of Sciences, Belorussian S.S.R., USSR
SOURCE: U.S.S.R. From: Otkrytiya, Izobret. 1991, (47), 20. CODEN: URXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Russian
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
SU 1699490	A1	19911223	SU 1989-4720902	19890720

PRIORITY APPLN. INFO.: SU 1989-4720902 19890720

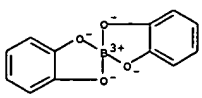
AB The title composition, conferring increased elasticity, radiation and heat resistant on membrane filters, comprises adipic acid-ε-caprolactam-hexamethylenediamine copolymer, glycerol, 2-aminopyridiniumdi(o-oxyphenylene) borate, and HCO2H.

IT 144672-73-1
RL: USES (Uses)
(polyamide composition containing, for membrane filters, for increased heat and radiation resistance)

RN 144672-73-1 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H^+

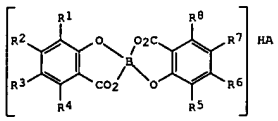
CM 2

CRN 504-29-0
CMF C5 H6 N2

L12 ANSWER 75 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1992:642782 CAPLUS
DOCUMENT NUMBER: 117:242782
TITLE: Thermal recording materials using disalicylborate double salt color developer
INVENTOR(S): Tanaka, Ginnosuke; Tottori, Kotaro; Shiraishi, Tomohisa; Kawai, Hajime
PATENT ASSIGNEE(S): Osaki Kogyo Co., Ltd., Japan; Yamada Kagaku Kogyo K. K.
SOURCE: Jpn. Kokai Tokkyo Koho, 8 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 04153082	A2	19920526	JP 1990-281310	19901018
PRIORITY APPLN. INFO.:			JP 1990-281310	19901018

GI



I

AB The title materials contain a leuco dye and, as a color developer, a double salt I m. 70-250° (R1,R4,R5,R8 = H, alkyl, aryl; R2,R3,R6,R7 = H, OH, alkyl, aryl; A = N-containing basic organic compound). A thermal recording paper using 2-anilino-3-methyl-6-ethylisopentylaminofluoran and I (R1-8 = H, A = Et2N) gave high-d. and low-fog images with good storage stability.

IT **69030-98-4**
RL: USES (Uses)
(color-developer, thermal recording material using)

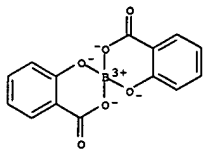
RN 69030-98-4 CAPLUS

CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS

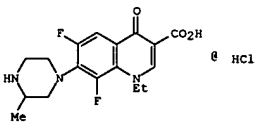
L12 ANSWER 75 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



CM 2

CRN 110-86-1
CMF C5 H5 N

L12 ANSWER 76 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1992:235582 CAPLUS
DOCUMENT NUMBER: 116:235582
TITLE: Improved preparation of lomefloxacin hydrochloride
AUTHOR(S): Wang, Erhua; Yao, Hong; Peng, Sixun
CORPORATE SOURCE: China Pharm. Univ., Nanjing, 210009, Peop. Rep. China
SOURCE: Zhongguo Yiyao Gongye Zazhi (1991), 22(10), 437-9
CODEN: ZYGZEA; ISSN: 1001-8255
DOCUMENT TYPE: Journal
LANGUAGE: Chinese
GI



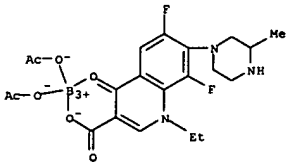
I

AB A facile preparation of lomefloxacin-HCl (I) starting from 2,3,4-trifluoronitrobenzene, via reduction, condensation-cyclization, ethylation, chelation, piperazination, hydrolysis, etc., was described. The reactions of all steps were carried out in very moderate conditions to afford overall yield of 40%.

IT **141301-51-1P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation and reaction with triethylamine)

RN 141301-51-1 CAPLUS

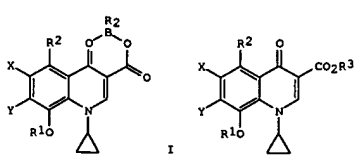
CN Boron, bis(acetato-O)[1-ethyl-6,8-difluoro-1,4-dihydro-7-(3-methyl-1-piperazinyl)-4-oxo-3-quinolinecarboxylato-O3,O4]-, (T-4)- (9CI) (CA INDEX NAME)



L12 ANSWER 77 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1992:152003 CAPLUS
DOCUMENT NUMBER: 116:152003
TITLE: (6,7-Substituted-8-alkoxy-1-cyclopropyl-1,4-dihydro-4-oxo-3-quinolinecarboxylic acid O3,O4)bis(acyloxy-O)borates and the salts thereof, and methods for their manufacture
INVENTOR(S): Takagi, Naomi; Fubasami, Hironobu; Matsukubo, Hiroshi
PATENT ASSIGNEE(S): Kyorin Pharmaceutical Co., Ltd., Japan
SOURCE: Eur. Pat. Appl., 13 pp.
CODEN: EPXXDW
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 464823	A1	19920108	EP 1991-111139	19910704
EP 464823	B1	19990922		
R: BE, CH, DE, ES, FR, GB, IT, LI, NL, SE				
JP 04069388	A2	19920304	JP 1990-178765	19900706
JP 07078065	B4	19950823		
US 5157117	A	19921020	US 1991-724164	19910701
ES 2137154	T3	19991216	ES 1991-111139	19910704
CA 2046361	AA	19920107	CA 1991-2046361	19910705
CA 2046361	C	19990720		
HU 58747	A2	19920330	HU 1991-2279	19910705
HU 215429	B	19990428		
AU 9180263	A1	19930128	AU 1991-80263	19910705
AU 646055	B2	19940203		
CN 1059527	A	19920318	CN 1991-104666	19910706
CN 1031795	B	19960515		
FI 103794	B1	19990930	FI 1992-12	19920102
AT 397656	B	19940627	AT 1992-9	19920107
AT 9200009	A	19931015		
PRIORITY APPLN. INFO.:			JP 1990-178765 A	19900706

OTHER SOURCE(S): CASREACT 116:152003; MARPAT 116:152003
GI



I II

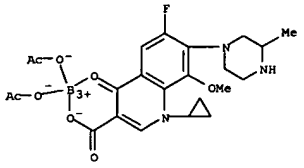
AB A process for the preparation of (6,7-substituted 8-alkoxy-1-cyclopropyl-1,4-

L12 ANSWER 77 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
dihydro-4-oxo-3-quinolinecarboxylic acid bis(acyloxy) borate of general formula I comprises the treatment of carboxyl compd. II with triacyloxyborate deriv. BR3. E.g., II (X = Y = F, R1 = Me, R2 = H, R3 = Et) (200 g) was added to a mixt. of boric acid (57.2 g) and zinc chloride (1.24 g) in acetic anhydride (300 mL) and acetic acid (400 mL) at a temp. of 50-60° to give 249 g of I (X = Y = F, R1 = Me, R2 = H, R = acetato).

IT **139693-53-1P 139693-54-2P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation and thermal reaction of)

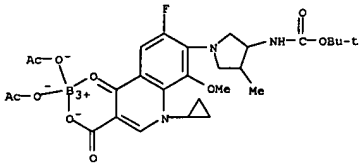
RN 139693-53-1 CAPLUS

CN Boron, bis(acetato-O)[1-cyclopropyl-6-fluoro-1,4-dihydro-8-methoxy-7-(3-methyl-1-piperazinyl)-4-oxo-3-quinolinecarboxylato-O3,O4]-, (T-4)- (9CI) (CA INDEX NAME)



RN 139693-54-2 CAPLUS

CN Boron, bis(acetato-O)[1-cyclopropyl-7-[3-[[(1,1-dimethylethoxy)carbonyl]amino]-4-methyl-1-pyrrolidinyl]-6-fluoro-1,4-dihydro-8-methoxy-4-oxo-3-quinolinecarboxylato-O3,O4]-, [T-4-(cis)]- (9CI)
(CA INDEX NAME)



L12 ANSWER 78 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1992:130450 CAPLUS

DOCUMENT NUMBER: 116:130450

TITLE: Spiroborate esters as fireproofing agents for plastics

INVENTOR(S): Horacek, Heinrich

PATENT ASSIGNEE(S): Chemie Linz (Deutschland) G.m.b.H., Germany

SOURCE: Ger. Offen., 6 pp.
CODEN: GWXXBX

DOCUMENT TYPE: Patent

LANGUAGE: German

FAMILY ACC. NUM. COUNT: 1

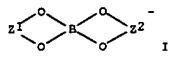
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
DE 4015490	A1	19911121	DE 1990-4015490	19900515

PRIORITY APPLN. INFO.: DE 1990-4015490 19900515

OTHER SOURCE(S): MARPAT 116:130450

GI



AB Amine salts of the spiroborates I [Z1, Z2 = diol residues or (jointly) a pentaerythritol residue] are halogen-free fireproofing agents for plastics. Thus, refluxing 0.24 mol H3BO3, 0.48 mol neopentyl glycol, and 360 mL EtOH for 2 h, adding 0.24 mol guanidine carbonate, and refluxing for 2 h gave neopentyl glycol spiroborate (2:1) guanidinium salt (1:1) (II). In a flame test (British Standard 5852, Crib 5), a polyurethane foam (bulk d. 0.021) containing 10% II and 10% melamine showed a weight loss of .apprx.2%; vs. 15% without II.

IT **139332-89-1P**, Neopentylglycol spiroborate (2:1) melamine salt (1:1) **139332-90-4P**, Manitol spiroborate (2:1) melamine salt (1:1) **139332-91-5P**, Pentaerythritol spiroborate (1:1) melamine salt (1:1) **139332-92-6P**, Pentaerythritol spiroborate (1:1) piperazine salt (1:1)
RL: PREP (Preparation)
(preparation of, as fireproofing agents for plastics)

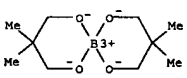
RN 139332-89-1 CAPLUS

CN Borate(1-), bis[2,2-dimethyl-1,3-propanediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 1,3,5-triazine-2,4,6-triamine (1:1) (9CI) (CA INDEX NAME)

CM 1

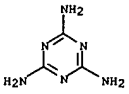
CRN 85283-52-9
CMF C10 H20 B O4 . H
CCI CCS

L12 ANSWER 78 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



CM 2

CRN 108-78-1
CMF C3 H6 N6

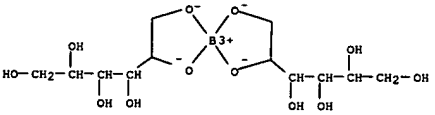


RN 139332-90-4 CAPLUS

CN Borate(1-), bis[D-mannitolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 1,3,5-triazine-2,4,6-triamine (1:1) (9CI) (CA INDEX NAME)

CM 1

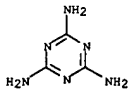
CRN 138662-85-8
CMF C12 H24 B O12 . H
CCI CCS



CM 2

CRN 108-78-1
CMF C3 H6 N6

L12 ANSWER 78 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

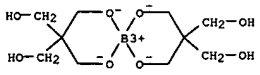


RN 139332-91-5 CAPLUS

CN Borate(1-), bis[2,2-bis(hydroxymethyl)-1,3-propanediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 1,3,5-triazine-2,4,6-triamine (1:1) (9CI) (CA INDEX NAME)

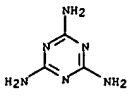
CM 1

CRN 138582-95-3
CMF C10 H20 B O8 . H
CCI CCS



CM 2

CRN 108-78-1
CMF C3 H6 N6



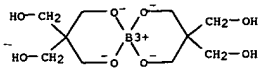
RN 139332-92-6 CAPLUS

CN Borate(1-), bis[2,2-bis(hydroxymethyl)-1,3-propanediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with piperazine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 138582-95-3
CMF C10 H20 B O8 . H
CCI CCS

L12 ANSWER 78 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



CM 2

CRN 110-85-0
CMF C4 H10 N2

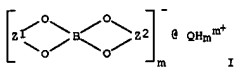
L12 ANSWER 79 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1992:85008 CAPLUS
DOCUMENT NUMBER: 116:85008
TITLE: Spirocyclic borate esters as flame retardants for plastics
INVENTOR(S): Horacek, Heinrich
PATENT ASSIGNEE(S): Chemie Linz G.m.b.H., Austria
SOURCE: Eur. Pat. Appl., 9 pp.
CODEN: EPXXDW
DOCUMENT TYPE: Patent
LANGUAGE: German
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 453821	A2	19911030	EP 1991-105131	19910330
EP 453821	A3	19920916		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE				
AT 9000945	A	19910915	AT 1990-945	19900424
AT 394370	B	19920325		
US 5147914	A	19920915	US 1991-678939	19910403
PL 168003	B1	19951230	PL 1991-289989	19910422
RO 110502	B1	19960130	RO 1991-147404	19910422
HU 57780	A2	19911230	HU 1991-1357	19910423
RU 2039764	C1	19950720	RU 1991-4895157	19910423

PRIORITY APPLN. INFO.: AT 1990-945 A 19900424

OTHER SOURCE(S): MARPAT 116:85008
GI



AB The spiroborates I [Q = N base; Z1, Z2 = residues of alcs. bearing 2-6 OH groups or together are C(CH2)4; m = 1-3] are halogen-free fireproofing agents for plastics. Refluxing 0.24 mol H3BO3, 0.48 g neopentyl glycol, and 360 mL EtOH for 2 h, cooling to 65°, slowly adding 0.12 mol guanidine carbonate, and refluxing for 2 h gave a spiroborate guanidinium salt (II). A polyurethane foam (1 kg) containing 10% II and 10% melamine, subjected to a burning test (BS 5852 Crib 5), lost 20 g, vs. 50 with 20 g melamine and no II.

IT **138582-97-5P 138582-98-6P 138582-99-7P**
138662-87-0P
RL: PREP (Preparation)
(manufacture of, as fireproofing agents for plastics)

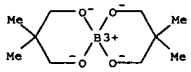
RN 138582-97-5 CAPLUS
CN Borate(1-), bis[2,2-dimethyl-1,3-propanediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 1,3,5-triazine-2,4,6-triamine (9CI) (CA INDEX NAME)

CM 1

CRN 85283-52-9

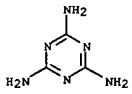
L12 ANSWER 79 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CMF C10 H20 B O4 . H
CCI CCS



CM 2

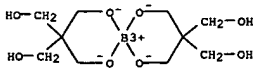
CRN 108-78-1
CMF C3 H6 N6



RN 138582-98-6 CAPLUS
CN Borate(1-), bis[2,2-bis(hydroxymethyl)-1,3-propanediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 1,3,5-triazine-2,4,6-triamine (9CI) (CA INDEX NAME)

CM 1

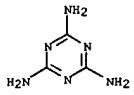
CRN 138582-95-3
CMF C10 H20 B O8 . H
CCI CCS



CM 2

CRN 108-78-1
CMF C3 H6 N6

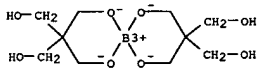
L12 ANSWER 79 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



RN 138582-99-7 CAPLUS
CN Borate(1-), bis[2,2-bis(hydroxymethyl)-1,3-propanediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with piperazine (9CI) (CA INDEX NAME)

CM 1

CRN 138582-95-3
CMF C10 H20 B O8 . H
CCI CCS



CM 2

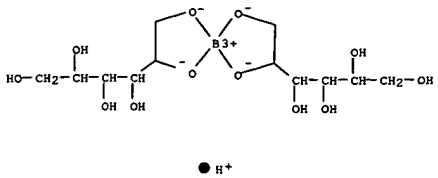
CRN 110-85-0
CMF C4 H10 N2

RN 138662-87-0 CAPLUS
CN Borate(1-), bis[D-mannitolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 1,3,5-triazine-2,4,6-triamine (9CI) (CA INDEX NAME)

CM 1

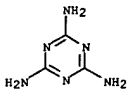
CRN 138662-85-8
CMF C12 H24 B O12 . H
CCI CCS

L12 ANSWER 79 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



CM 2

CRN 108-78-1
CMF C3 H6 N6



L12 ANSWER 80 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1992:5952 CAPLUS
DOCUMENT NUMBER: 116:5952
TITLE: NMR study of the effect of nitrogen-borane coordination on the conformational equilibrium of six membered ring heterocycles
AUTHOR(S): Flores-Parra, Angelina; Farfan, Norberto; Hernandez-Bautista, Alberto I.; Fernandez-Sanchez, Lilia; Contreras, Rosalinda
CORPORATE SOURCE: Cent. Invest. Estud. Avanzados, Inst. Politec. Nac., Mexico City, 07000, Mex.
SOURCE: Tetrahedron (1991), 47(34), 6903-14
CODEN: TETRAB; ISSN: 0040-4020
DOCUMENT TYPE: Journal
LANGUAGE: English

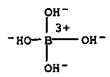
AB The syntheses, conformational and spectroscopic studies of N-borane adducts of 14 nitrogen-containing six-membered ring heterocycles are reported.
It was found that borane can act as a conformational and configurational locking agent. In addition, it can be very helpful for the assignment of the
chemical shifts of other atoms or groups in the mol. as well as to ascertain
the configuration at substituted carbons.

IT **137546-85-1P**
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(preparation and NMR of, as analog of borane coordination compound)

RN 137546-85-1 CAPLUS
CN Borate(1-), tetrahydroxy-, hydrogen, compd. with 1,2,2,6,6-pentamethylpiperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

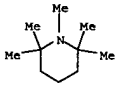
CRN 137546-84-0
CMF B H4 O4 . H
CCI CCS



CM 2

CRN 79-55-0
CMF C10 H21 N

L12 ANSWER 80 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

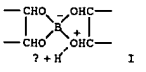


L12 ANSWER 81 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1991:472980 CAPLUS
DOCUMENT NUMBER: 115:72980
TITLE: Surface modification of inorganic fireproofing agents
INVENTOR(S): Hamanaka, Hiroyoshi
PATENT ASSIGNEE(S): Boron International K. K., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 15 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 02294389	A2	19901205	JP 1989-113656	19890508

PRIORITY APPLN. INFO.: JP 1989-113656 19890508

GI



I

AB Fireproofing agents for polymers with good compatibility are prepared by forming on hydrated metal compds. hybrid films of oriented, water-insol. hydrocarbyl compds. and semipolar B compds. containing I groups and/or basic N
compound complexes. Heating 1:1 di(glycerol)borate and ricinoleic acid dimer at 230-240° for 6 h, cooling to 70°, adding MEK and ethoxylated dihexadecylamine, and heating at 70-75° for 1 and at 120-130°/150 mm for 2 h gave a brown, viscous complex. Mixing 40 g 25% tricetylaluminum solution and 20 g 50% solution of this complexwith 1 kg
Al(OH)3 particles (1 μm) and stripping solvents gave modified Al(OH)3, which was kneaded g0:40 with polyethylene to give a 1-mm sheet with good acid and flex resistance.

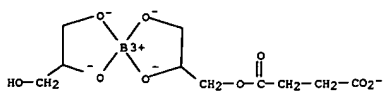
IT **134620-04-5**
RL: USES (Uses)
(surface modification by, of inorg. fireproofing agents)

RN 134620-04-5 CAPLUS
CN Borate(2-), [mono(2,3-dihydroxypropyl) butanedioato(3-)][1,2,3-propanetriolato(2-)-O1,O2]-, (T-4)-, dihydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 106530-70-5
CMF C10 H15 B O9 . 2 H
CCI CCS

L12 ANSWER 81 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



●2 H+

CM 2

CRN 110-86-1
CMF C5 H5 N

L12 ANSWER 82 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1991:115064 CAPLUS
DOCUMENT NUMBER: 114:115064
TITLE: Preparation and use of polyoxometallates for treatment of retrovirus infections
INVENTOR(S): Hill, Craig L.; Schinazi, Raymond F.
PATENT ASSIGNEE(S): Johnson Matthey PLC, UK
SOURCE: Eur. Pat. Appl., 15 pp.
CODEN: EPXXDW
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 360619	A2	19900328	EP 1989-309678	19890922
EP 360619	A3	19910807		
R: AT, BE, CH, DE, ES, FR, GB, GR, IT, LI, LU, NL, SE				
WO 9003176	A1	19900405	WO 1989-US4028	19890922
W: AU, DK, FI, HU, JP, KR, NO				
AU 8943168	A1	19900418	AU 1989-43168	19890922
ZA 8907260	A	19900725	ZA 1989-7260	19890922
US 6911470	B1	20050628	US 1993-140885	19931025
PRIORITY APPLN. INFO.:			US 1988-247641 A	19880922
			WO 1989-US4028 A	19890922
			US 1990-474389 B1	19900205
			US 1993-24837 B1	19930301

OTHER SOURCE(S): MARPAT 114:115064

AB The title compds. comprise e.g. (X+)n(H+)4-n(SiW12O40)4- (n = 0-4; X = K, Na, NH4, histidinium, lysinium, argininium), (X+)n(H+)5-m(BW12O40)5- (m = 0-5; X as above), (X+)n(H+)4-n(W10O32)4- (n = 0-4; X as above), (X+)p(H+)3-p(PW12O40)3- (p = 0-3; X as above), and their hydrates. X can generally be any naturally occurring monocationic amino acid or polypeptide containing arginine, histidine, or lysine. Thus, water-saturated NaSiO3 and Na2WO4.2H2O were reacted to prepare H4SiW12O40, which was then treated with lysine-HCl to form the corresponding lysinium salt. Antiviral activities of the compds. of the invention, in a variety of assays (in vitro and in vivo), are tabulated. Cytotoxicity and bone marrow toxicity of the compds. were determined The compds. are useful against e.g. human immunodeficiency virus infection.

IT **131541-70-3P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of, for retrovirus infection treatment)

RN 131541-70-3 CAPLUS

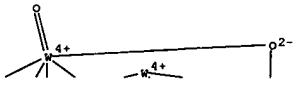
CN L-Histidine, tetracosa-μ-oxododecaoxo[μ12-[tetrahydroxyborato(5-)-O:O:O:O':O':O':O'':O'':O'':O''':O''':O''']]dodecatungstate(5-) (9CI) (CA INDEX NAME)

CM 1

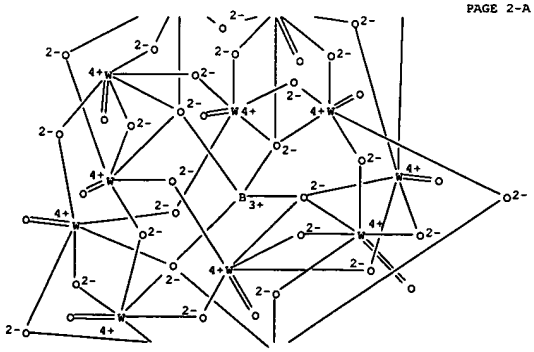
CRN 12297-12-0
CMF B O40 W12 . 5 H
CCI CCS

L12 ANSWER 82 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

PAGE 1-A

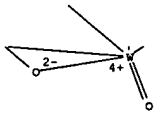


PAGE 2-A



L12 ANSWER 82 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

PAGE 3-A

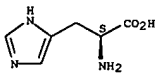


●5 H+

CM 2

CRN 71-00-1
CMF C6 H9 N3 O2

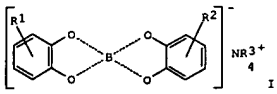
Absolute stereochemistry. Rotation (-).



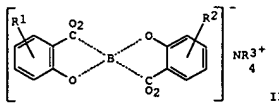
L12 ANSWER 83 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1991:93563 CAPLUS
DOCUMENT NUMBER: 114:93563
TITLE: Supporting electrolyte composition for manufacture of solid electrolytic capacitor
INVENTOR(S): Koseki, Tetsuya; Tsuji, Tatsunori; Yokoyama, Yutaka
PATENT ASSIGNEE(S): Nippon Chemi-Con Corp., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 4 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 02166713	A2	19900627	JP 1988-322435	19881221
JP 3076872	B2	20000814		
PRIORITY APPLN. INFO.:			JP 1988-322435	19881221

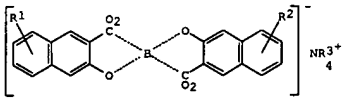
OTHER SOURCE(S): MARPAT 114:93563
GI



I



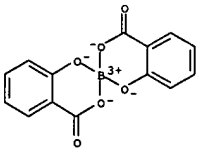
II



III

AB The title composition, which is used in electrolytic polymerization for preparing a solid electrolytic layer on a valve metal oxide layer, comprises an aromatic hydroxy compound B complex quaternary ammonium salt I or an aromatic hydroxycarboxylic acid B complex quaternary ammonium salt II or III (R1-2 = H, alkyl, OH, amino, CO2H; R3 = H, C1-5 alkyl; R2-3 may form a cyclic ammonium). An Al foil was coated with a dielec. oxide layer and used as an anode in electrolytic polymerization of pyrrole in the presence of N,N-dimethylpiperidinium borodisalicylate to give a solid electrolyte. A solid electrolytic capacitor using the electrolyte showed thermally stable

L12 ANSWER 83 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



2 (D1—Me)

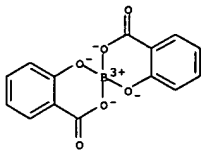
CM 2

CRN 15312-12-6
CMF C6 H14 N

L12 ANSWER 83 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
elec. characteristics.
IT 131121-99-8 132099-07-1
RL: TEM (Technical or engineered material use); USES (Uses)
(supporting electrolyte, for electrolytic polymerization, for solid electrolytic capacitor)
RN 131121-99-8 CAPLUS
CN Piperidinium, 1,1-dimethyl-, (T-4)-bis[2-hydroxybenzoato(2-)-O1,O2]borate(1-) (9CI) (CA INDEX NAME)

CM 1

CRN 38403-08-6
CMF C14 H8 B O6
CCI CCS



CM 2

CRN 15302-91-7
CMF C7 H16 N

RN 132099-07-1 CAPLUS
CN Pyrrolidinium, 1,1-dimethyl-, bis[2-hydroxymethylbenzoato(2-)-O1,O2]borate(1-) (9CI) (CA INDEX NAME)

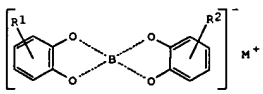
CM 1

CRN 132099-06-0
CMF C16 H12 B O6
CCI CCS, IDS

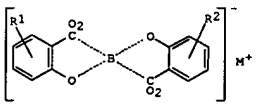
L12 ANSWER 84 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1991:93562 CAPLUS
DOCUMENT NUMBER: 114:93562
TITLE: Supporting electrolyte composition for manufacture of solid electrolytic capacitor
INVENTOR(S): Koseki, Tetsuya; Tsuji, Tatsunori; Yokoyama, Yutaka
PATENT ASSIGNEE(S): Nippon Chemi-Con Corp., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 4 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 02166712	A2	19900627	JP 1988-322436	19881221
JP 3076873	B2	20000814		
PRIORITY APPLN. INFO.:			JP 1988-322436	19881221

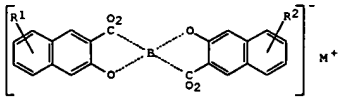
OTHER SOURCE(S): MARPAT 114:93562
GI



I



II



III

AB The title composition, which is used in electrolytic polymerization for preparing a solid electrolytic layer on a valve metal oxide layer, comprises an aromatic hydroxy compound B complex amine salt I or an aromatic hydroxycarboxylic acid B complex amine salt II or III (R1-2 = H, alkyl, OH, amino, CO2H; M = amine). An Al foil was coated with a dielec. oxide layer and used as an anode in electrolytic polymerization of pyrrole in the presence of Et3N borodiresorcylate to give a solid electrolyte. A solid electrolytic capacitor using the electrolyte showed thermally stable elec. characteristics.

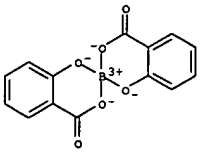
L12 ANSWER 84 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

IT **131122-00-4 131122-01-5**
RL: TEM (Technical or engineered material use); USES (Uses)
(supporting electrolyte, for electrolytic polymerization, for solid electrolytic capacitor)

RN 131122-00-4 CAPLUS
CN Borate(1-), bis[2-hydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 1-ethylpiperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS



● H+

CM 2

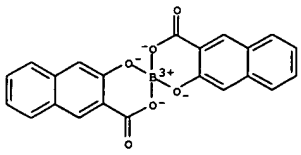
CRN 766-09-6
CMF C7 H15 N

RN 131122-01-5 CAPLUS
CN Pyrrolidinium, 1,1-dimethyl-, (T-4)-bis[3-hydroxy-2-naphthalenecarboxylato(2-)-O2,O3]borate(1-) (9CI) (CA INDEX NAME)

CM 1

CRN 47611-21-2
CMF C22 H12 B O6
CCI CCS

L12 ANSWER 84 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



CM 2

CRN 15312-12-6
CMF C6 H14 N

L12 ANSWER 85 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1991:43249 CAPLUS
DOCUMENT NUMBER: 114:43249
TITLE: Compounds of borotungstic acid with nitrogen compounds
AUTHOR(S): Chaigneau, M.
CORPORATE SOURCE: Lab. Gaz, Paris, F 75270, Fr.
SOURCE: Annales Pharmaceutiques Francaises (1990), 48(1), 42-4
CODEN: APFRAD; ISSN: 0003-4509
DOCUMENT TYPE: Journal
LANGUAGE: French

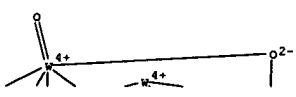
AB Alkaloids such as quinine, cinchonine, brucine, morphine, strychnine, and nitrogenous compds., e.g., procaine, choline, were allowed to react with borotungstic acid and shown to form borotungstates.

IT **130555-54-3P 130555-55-4P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)

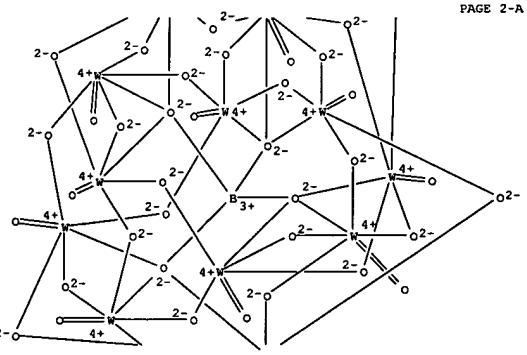
RN 130555-54-3 CAPLUS
CN Tungstate(5-), tetracosa-μ-oxododecaoxo[μ12-[tetrahydroxyborato(5-)-O:O:O:O':O':O':O'':O'':O'':O''':O''':O''']]dodeca-, pentahydrogen, compd. with (3S-cis)-3-ethyldihydro-4-[(1-methyl-1H-imidazol-5-yl)methyl]-2(3H)-furanone (1:5) (9CI) (CA INDEX NAME)

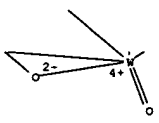
CM 1

CRN 12297-12-0
CMF B O40 W12 . 5 H
CCI CCS



L12 ANSWER 85 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



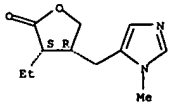


●5 H+

CM 2

CRN 92-13-7
CMF C11 H16 N2 O2

Absolute stereochemistry.

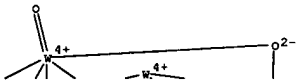


RN 130555-55-4 CAPLUS
CN Tungstate(5-), tetracosa-μ-oxododecaoxo[μ12-[tetrahydroxyborato(5-)-O:O:O:O':O':O'':O'':O'':O''':O''':O''']]dodeca-, pentahydrogen, compd. with (S)-3-(1-methyl-2-pyrrolidinyl)pyridine (1:4) (9CI) (CA INDEX NAME)

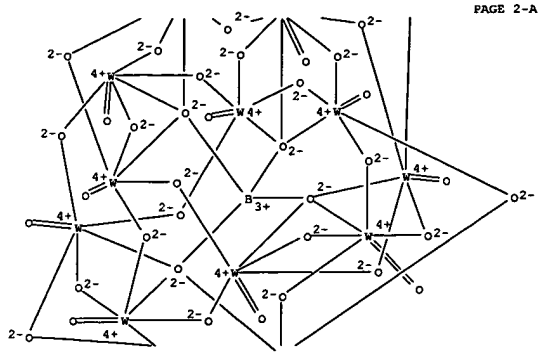
CM 1

CRN 12297-12-0
CMF B O40 W12 . 5 H
CCI CCS

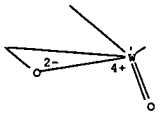
PAGE 1-A



PAGE 2-A



PAGE 3-A

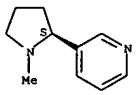


●5 H+

CM 2

CRN 54-11-5
CMF C10 H14 N2

Absolute stereochemistry. Rotation (-).



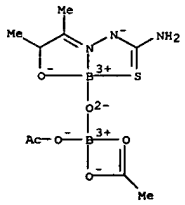
L12 ANSWER 86 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1990:244848 CAPLUS
DOCUMENT NUMBER: 112:244848
TITLE: Synthetic and structural studies of some tetracoordinated boron complexes of bifunctional tridentate Schiff bases
AUTHOR(S): Singh, V. P.; Singh, R. V.; Tandon, J. P.
CORPORATE SOURCE: Dep. Chem., Univ. Rajasthan, Jaipur, 302004, India
SOURCE: Journal fuer Praktische Chemie (Leipzig) (1989), 331(4), 690-6
CODEN: JPCEAO; ISSN: 0021-8383
DOCUMENT TYPE: Journal
LANGUAGE: English

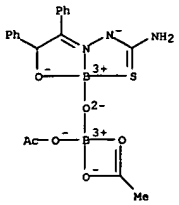
AB [(AcO)2B]2O reacted with RR1C:NNHC(S)NH2 (H2L; R = CHOHCH3, R1 = Me; R = 2-C6H4OH, R1 = H, Me; R = 2-hydroxy-1-naphthyl, R1 = H; R = CHOHPh, R1 = Ph) to give (AcO)2BOBL and [LB]2O. L are tridentate, coordinating through the O, S, and N atoms. The complexes were characterized by IR and 11B and 1H NMR spectra.

IT **127465-88-7P 127465-92-3P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)

RN 127465-88-7 CAPLUS
CN Boron, (acetato-O)(acetato-O,O')[2-(2-hydroxy-1-methylpropylidene)hydrazinecarbothioamidato(2-)-N2,O2,S]-μ-oxodi- (9CI) (CA INDEX NAME)



RN 127465-92-3 CAPLUS
CN Boron, (acetato-O)(acetato-O,O')[2-(2-hydroxy-1,2-diphenylethylidene)hydrazinecarbothioamidato(2-)-N2,O2,S]-μ-oxodi- (9CI) (CA INDEX NAME)

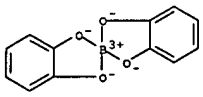


ACCESSION NUMBER: 1989:632934 CAPLUS
DOCUMENT NUMBER: 111:232934
TITLE: Studies on spiroborate complexes. Part III. Fast-atom-bombardment mass spectrometry of bis-catechol spiroborate and its analogs
AUTHOR(S): Okamoto, Yoshihisa; Takei, Yuka; Rose, Malcolm E.
CORPORATE SOURCE: Coll. Lib. Arts Sci., Kitasato Univ., Sagamihara, 228, Japan
SOURCE: International Journal of Mass Spectrometry and Ion Processes (1989), 87(2), 225-35
CODEN: IJMPDN; ISSN: 0168-1176
DOCUMENT TYPE: Journal
LANGUAGE: English
AB Neg.-ion fast-atom-bombardment mass spectra of 40 spiroborate complexes are simple and highly diagnostic of the structures of the borate anions. In hydroxylic liquid matrixes, exchange of the spiroborate ligand and solvent mols. occurs slowly on the probe tip. This interference is particularly noticeable with glycerol as solvent. The repercussions of in situ ligand exchange for studies of other anions are discussed. Pos.-ion fast-atom-bombardment mass spectra can be used to identify readily the counter-cations in the spiroborate complexes.
IT **111932-23-1 111932-24-2 111932-28-6 111932-30-0 111932-31-1 111932-32-2 111932-34-4 111932-35-5 123738-44-3 123738-46-5 123738-48-7 123778-58-5 123778-60-9**
RL: PRP (Properties)
(fast-atom-bombardment mass spectrum of)

RN 111932-23-1 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H+

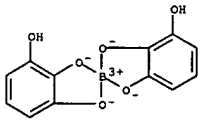
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-24-2 CAPLUS
CN Borate(1-), bis[1,2,3-benzenetriolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 46944-61-0
CMF C12 H8 B O6 . H
CCI CCS



● H+

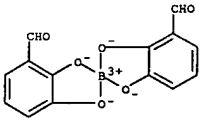
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-28-6 CAPLUS
CN Borate(1-), bis[2,3-dihydroxybenzaldehydato(2-)-O2,O3]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 111932-27-5
CMF C14 H8 B O6 . H
CCI CCS



● H+

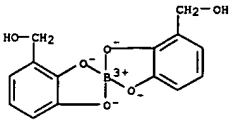
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-30-0 CAPLUS
CN Borate(1-), bis[3-(hydroxymethyl)-1,2-benzenediolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 111932-29-7
CMF C14 H12 B O6 . H
CCI CCS



● H+

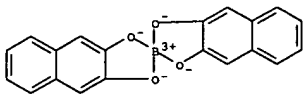
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-31-1 CAPLUS
CN Borate(1-), bis[2,3-naphthalenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 47422-29-7
CMF C20 H12 B O4 . H
CCI CCS



● H+

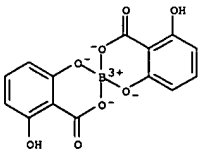
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-32-2 CAPLUS
CN Borate(1-), bis[2-hydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1



● H+

CM 2

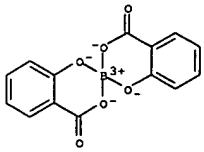
CRN 695-34-1
CMF C6 H8 N2

RN 111932-35-5 CAPLUS
CN Borate(1-), bis[[1,1'-biphenyl]-2,2'-diolato(2-)-O,O']-, (T-4)-, hydrogen,
compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 53993-02-5
CMF C24 H16 B O4 . H
CCI CCS

CMF C14 H8 B O6 . H
CCI CCS



● H+

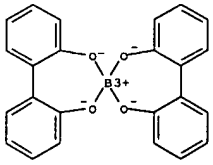
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-34-4 CAPLUS
CN Borate(1-), bis[2,6-dihydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 95692-94-7
CMF C14 H8 B O8 . H
CCI CCS



● H+

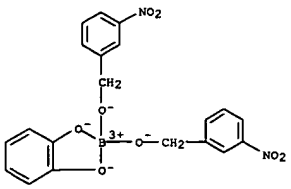
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 123738-44-3 CAPLUS
CN Borate(1-), [1,2-benzenediolato(2-)-O,O']bis(3-nitrobenzenemethanolato-Oα)-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 123738-43-2
CMF C20 H16 B N2 O8 . H
CCI CCS



● H^+

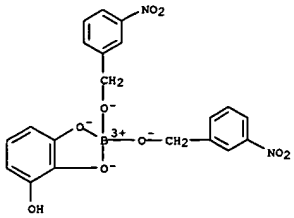
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 123738-46-5 CAPLUS
CN Borate(1-), {1,2,3-benzenetriolato(2-)-O1,O2]bis(3-nitrobenzenemethanolato-Oα)-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 123738-45-4
CMF C20 H16 B N2 O9 . H
CCI CCS



● H^+

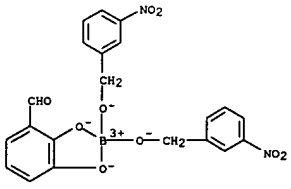
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 123738-48-7 CAPLUS
CN Borate(1-), (2,3-dihydroxybenzaldehydato(2-)-O2,O3]bis(3-nitrobenzenemethanolato-Oα)-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 123738-47-6
CMF C21 H16 B N2 O9 . H
CCI CCS



● H^+

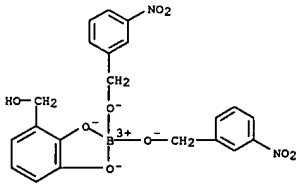
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 123778-58-5 CAPLUS
CN Borate(1-), [3-(hydroxymethyl)-1,2-benzenediolato(2-)-O1,O2]bis(3-nitrobenzenemethanolato-Oα)-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 123778-57-4
CMF C21 H18 B N2 O9 . H
CCI CCS



● H^+

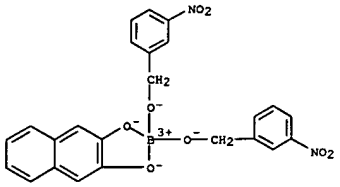
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 123778-60-9 CAPLUS
CN Borate(1-), [2,3-naphthalenediolato(2-)-O,O']bis(3-nitrobenzenemethanolato-Oα)-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 123778-59-6
CMF C24 H18 B N2 O8 . H
CCI CCS



● H+

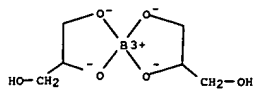
CM 2

CRN 695-34-1
CMF C6 H8 N2

IT **123738-49-8P 123778-62-1P 123778-64-3P**
123778-66-5P
RL: PRP (Properties); PREP (Preparation)
(formation and fast-atom-bombardment mass spectrum of)
RN 123738-49-8 CAPLUS
CN Borate(1-), bis[1,2,3-propanetriolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 49625-59-4
CMF C6 H12 B O6 . H
CCI CCS



● H+

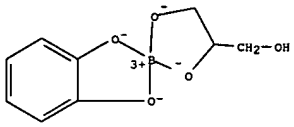
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 123778-62-1 CAPLUS
CN Borate(1-), [1,2-benzenediolato(2-)-O,O'][1,2,3-propanetriolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 123778-61-0
CMF C9 H10 B O5 . H
CCI CCS



● H+

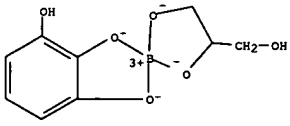
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 123778-64-3 CAPLUS
CN Borate(1-), [1,2,3-benzenetriolato(2-)-O1,O2][1,2,3-propanetriolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI)
(CA INDEX NAME)

CM 1

CRN 123778-63-2
CMF C9 H10 B O6 . H
CCI CCS



● H+

CM 2

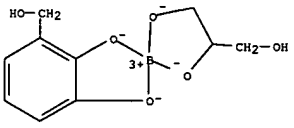
CRN 695-34-1
CMF C6 H8 N2

RN 123778-66-5 CAPLUS
CN Borate(1-), [3-(hydroxymethyl)-1,2-benzenediolato(2-)-O1,O2][1,2,3-propanetriolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with

4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 123778-65-4
CMF C10 H12 B O6 . H
CCI CCS



● H+

CM 2

CRN 695-34-1
CMF C6 H8 N2

L12 ANSWER 88 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1989:96437 CAPLUS
DOCUMENT NUMBER: 110:96437
TITLE: Organoboron antioxidants. Part 1. Boric acid derivatives as primary antioxidants
AUTHOR(S): Koenig, T.; Maennel, D.; Habicher, W. D.; Schwetlick, K.
CORPORATE SOURCE: Dep. Chem., Dresden Univ. Technol., Dresden, DDR-8027,
Ger. Dem. Rep.
SOURCE: Polymer Degradation and Stability (1988), 22(2), 137-45
CODEN: PDSTDW; ISSN: 0141-3910
DOCUMENT TYPE: Journal
LANGUAGE: English

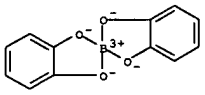
AB Organoboron compds., particularly boric acid esters and arylboronic acid derivs., are known to be good antioxidants. The inhibiting activities of various B compds. are compared with those of phenolic antioxidants and phosphites. Whereas in the autoxidn. of polypropylene, aryl borates give rise to approx. the same induction periods as the constituent phenols, aryl phenylboronates are, to some extent, considerably better. For the calcn. of kinetic data, particularly the rate consts. (k7) of the reaction with peroxyl radicals and the stoichiometric factor (f), investigations into the inhibiting activity of B compds. in the initiated oxidation of cumene at low temps. are used. The rate constant k7 of aryl phenylboronates is approx. 10 times greater than that for structurally analogous phosphites, whereas aryl borates can be regarded as depots for phenols.

IT **119177-56-9**
RL: USES (Uses)
(antioxidants, for cumene and polypropylene)

RN 119177-56-9 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 2,2,6,6-tetramethyl-4-piperidinone (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS

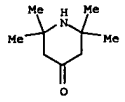


● H^+

CM 2

CRN 826-36-8
CMF C9 H17 N O

L12 ANSWER 88 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



L12 ANSWER 89 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1988:406562 CAPLUS
DOCUMENT NUMBER: 109:6562
TITLE: Reactions of boric acid with 3-formylsalicylic acid, glycolic acid, and benzilic acid leading to the synthesis of 3-formylsalicylato-, glycolato-, and benzilato-borate (III) salts
AUTHOR(S): Dey, K.; Gangopadhyay, A.; Biswas, A. K.
CORPORATE SOURCE: Dep. Chem., Univ. Kalyani, Kalyani, 741 235, India
SOURCE: Journal of Bangladesh Academy of Sciences (1987), 11(1), 55-66
CODEN: JBACDF; ISSN: 0378-8121
DOCUMENT TYPE: Journal
LANGUAGE: English

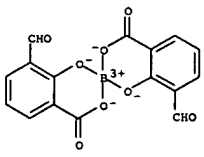
AB Reaction of B(OH)3 with 3-formylsalicylic acid, glycolic acid, and benzilic acid in the presence of various large cations afforded bis- and mono-3-formylsalicylato-, glycolato- and benzilatoborate(III) salts with large cations. The phys. and spectroscopic properties of these new borate(III) salts are reported. Synthesis and characterization of the neutral four-coordinate boron compound [LB(OH)(py)] (L = 3-formylsalicylato) is also discussed.

IT **114803-05-3P 114803-10-0P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)

RN 114803-05-3 CAPLUS
CN Borate(1-), bis[3-formyl-2-hydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 114803-04-2
CMF C16 H8 B O8 . H
CCI CCS



● H^+

CM 2

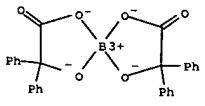
CRN 110-86-1
CMF C5 H5 N

L12 ANSWER 89 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

RN 114803-10-0 CAPLUS
CN Borate(1-), bis[α-hydroxy-α-phenylbenzeneacetato(2-)]-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 95692-86-7
CMF C28 H20 B O6 . H
CCI CCS



● H^+

CM 2

CRN 110-86-1
CMF C5 H5 N

L12 ANSWER 90 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1988:84752 CAPLUS
DOCUMENT NUMBER: 108:84752
TITLE: Studies on spiroborate complexes - II. Structural elucidation of biscatechol spiroborate and its analogs by NMR and mass spectral spectroscopy
AUTHOR(S): Okamoto, Yoshihisa; Takei, Yuka; Takagi, Kaname
CORPORATE SOURCE: Coll. Lib. Arts Sci., Kitasato Univ., Sagamihara, 228, Japan
SOURCE: Polyhedron (1987), 6(12), 2119-28
CODEN: PLYHDE; ISSN: 0277-5387
DOCUMENT TYPE: Journal
LANGUAGE: English

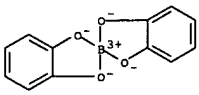
AB Nine spiroborate complexes were analyzed by NMR and mass spectroscopy, and the characteristic spectra and fragmentation patterns are reported. On the basis of these results, the structure of J. Meulenhoff's (1925) free acid was reinvestigated.

IT **111932-23-1 111932-24-2 111932-26-4 111932-28-6 111932-30-0 111932-31-1 111932-32-2 111932-34-4 111932-35-5**
RL: PRP (Properties)
(NMR and mass spectra and structure of)

RN 111932-23-1 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS

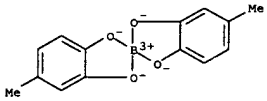


● H^+

CM 2

CRN 695-34-1
CMF C6 H8 N2

L12 ANSWER 90 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



● H^+

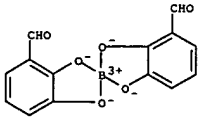
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-28-6 CAPLUS
CN Borate(1-), bis[2,3-dihydroxybenzaldehydato(2-)-O2,O3]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 111932-27-5
CMF C14 H8 B O6 . H
CCI CCS



● H^+

CM 2

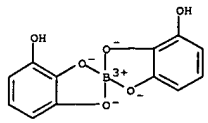
CRN 695-34-1

L12 ANSWER 90 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

RN 111932-24-2 CAPLUS
CN Borate(1-), bis[1,2,3-benzenetriolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 46944-61-0
CMF C12 H8 B O6 . H
CCI CCS



● H^+

CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-26-4 CAPLUS
CN Borate(1-), bis[4-methyl-1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

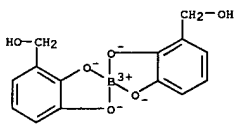
CRN 111932-25-3
CMF C14 H12 B O4 . H
CCI CCS

L12 ANSWER 90 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
CMF C6 H8 N2

RN 111932-30-0 CAPLUS
CN Borate(1-), bis[3-(hydroxymethyl)-1,2-benzenediolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 111932-29-7
CMF C14 H12 B O6 . H
CCI CCS



● H^+

CM 2

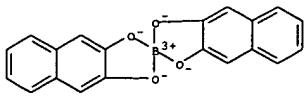
CRN 695-34-1
CMF C6 H8 N2

RN 111932-31-1 CAPLUS
CN Borate(1-), bis[2,3-naphthalenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 47422-29-7

L12 ANSWER 90 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
CMF C20 H12 B O4 . H
CCI CCS



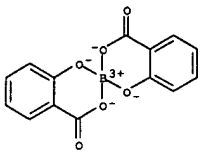
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-32-2 CAPLUS
CN Borate(1-), bis[2-hydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

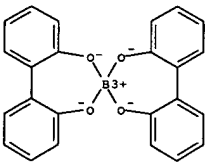
CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS



L12 ANSWER 90 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 53993-02-5
CMF C24 H16 B O4 . H
CCI CCS



CM 2

CRN 695-34-1
CMF C6 H8 N2

L12 ANSWER 90 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

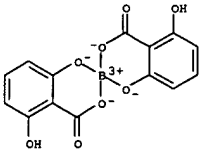
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-34-4 CAPLUS
CN Borate(1-), bis[2,6-dihydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 95692-94-7
CMF C14 H8 B O8 . H
CCI CCS

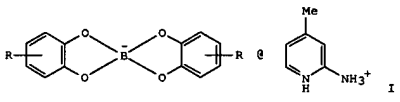


CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-35-5 CAPLUS
CN Borate(1-), bis[[1,1'-biphenyl]-2,2'-diolato(2-)-O,O']-, (T-4)-, hydrogen,

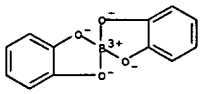
L12 ANSWER 91 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1988:21943 CAPLUS
DOCUMENT NUMBER: 108:21943
TITLE: Studies on spiroborate complexes - I. A new synthesis
of bis-catechol spiroborate and its analogs using 2-amino-4-methylpyridine borane
AUTHOR(S): Okamoto, Yoshihisa; Kinoshita, Toshio; Takei, Yuka; Matsumoto, Yoshio
CORPORATE SOURCE: Sch. Pharm. Sci., Kitasato Univ., Tokyo, 108, Japan
SOURCE: Polyhedron (1986), 5(12), 2051-7
CODEN: PLYHDE; ISSN: 0277-5387
DOCUMENT TYPE: Journal
LANGUAGE: English
GI



AB A new synthesis of bis-catechol spiroborate and its analogs are described.
Five-, six- and seven-membered spiroborates are obtained. Thus, reaction of 1,2-(HO)2C6H3R (R = H, 3-OH, 4-Me, 3-CHO) with 2-amino-4-methylpyridine
borane in HC(OEt)3 yielded spiroborates I (same R). Spiroborates of 2,3-dihydroxynaphthalene, salicylic acid, 2,6-dihydroxybenzoic acid, and 2,2'-dihydroxybiphenyl were also prepared A mechanism is proposed where the
reaction is initiated by removal of a phenolic proton by the orthoformate solvent.
IT **111932-23-1P 111932-24-2P 111932-26-4P**
111932-28-6P 111932-30-0P 111932-31-1P
111932-32-2P 111932-34-4P 111932-35-5P
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 111932-23-1 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H+

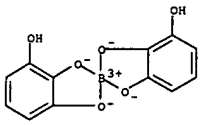
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-24-2 CAPLUS
CN Borate(1-), bis[1,2,3-benzenetriolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

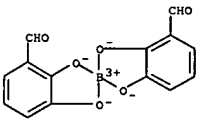
CRN 46944-61-0
CMF C12 H8 B O6 . H
CCI CCS



● H+

CM 2

CRN 695-34-1
CMF C6 H8 N2



● H+

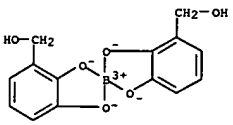
CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-30-0 CAPLUS
CN Borate(1-), bis[3-(hydroxymethyl)-1,2-benzenediolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 111932-29-7
CMF C14 H12 B O6 . H
CCI CCS



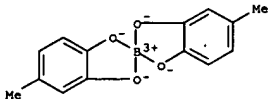
● H+

CM 2

RN 111932-26-4 CAPLUS
CN Borate(1-), bis[4-methyl-1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 111932-25-3
CMF C14 H12 B O4 . H
CCI CCS



● H+

CM 2

CRN 695-34-1
CMF C6 H8 N2

RN 111932-28-6 CAPLUS
CN Borate(1-), bis[2,3-dihydroxybenzaldehydato(2-)-O2,O3]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

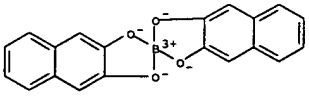
CRN 111932-27-5
CMF C14 H8 B O6 . H
CCI CCS

CRN 695-34-1
CMF C6 H8 N2

RN 111932-31-1 CAPLUS
CN Borate(1-), bis[2,3-naphthalenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 47422-29-7
CMF C20 H12 B O4 . H
CCI CCS



● H+

CM 2

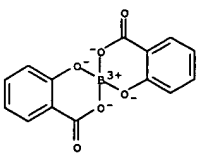
CRN 695-34-1
CMF C6 H8 N2

RN 111932-32-2 CAPLUS
CN Borate(1-), bis[2-hydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H

L12 ANSWER 91 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
CCI CCS



CM 2

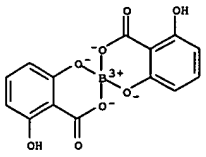
CRN 695-34-1
CMF C6 H8 N2

RN 111932-34-4 CAPLUS
CN Borate(1-), bis[2,6-dihydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 95692-94-7
CMF C14 H8 B O8 . H
CCI CCS

L12 ANSWER 91 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



CM 2

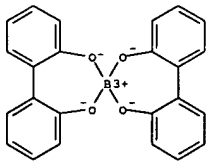
CRN 695-34-1
CMF C6 H8 N2

RN 111932-35-5 CAPLUS
CN Borate(1-), bis[[1,1'-biphenyl]-2,2'-diolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 4-methyl-2-pyridinamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 53993-02-5
CMF C24 H16 B O4 . H
CCI CCS

L12 ANSWER 91 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



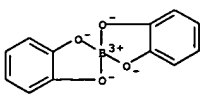
CM 2

CRN 695-34-1
CMF C6 H8 N2

L12 ANSWER 92 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1987:206657 CAPLUS
DOCUMENT NUMBER: 106:206657
TITLE: Synthesis and study of new boron complexes
AUTHOR(S): Grachek, V. I.; Naumova, S. F.; Shelemova, I. V.
CORPORATE SOURCE: Inst. Fiz.-Org. Khim., Minsk, USSR
SOURCE: Vestsi Akademii Navuk BSSR, Seryya Khimichnykh Navuk (1987), (1), 116-18
CODEN: VBSKAK; ISSN: 0002-3590
DOCUMENT TYPE: Journal
LANGUAGE: Russian
AB H3BO3 reacted with pyrocatechol (H2L) in presence of N bases to give HQ[BL2] (Q = 2,4-lutidine, 2-methylquinoline, piperazine, nicotinamide, 2-hydroxypyrrolidine). The complexes were characterized by IR spectra. The herbicidal and antibacterial activities of HQ[BL2] and the quinoline, piperidine and pyridine analogs were studied. Compds. containing a pyridine ring are more effective antibacterial agents; the quinoline and piperidene derivs. are the more effective herbicides.
IT **108339-50-0 108339-52-2**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); BIOL (Biological study)
(antibacterial and herbicidal activity of)
RN 108339-50-0 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with piperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



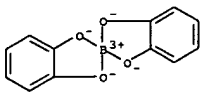
CM 2

CRN 110-89-4
CMF C5 H11 N

RN 108339-52-2 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H^+

CM 2

CRN 110-86-1
CMF C5 H5 N

IT **108339-53-3P 108339-55-5P 108339-56-6P 108357-04-6P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation and antibacterial and herbicidal activity of)

RN 108339-53-3 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with 2,4-dimethylpyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS

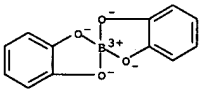
RN 108339-56-6 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 2-pyrrolidinol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 67318-87-0
CMF C4 H9 N O

CM 2

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS

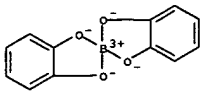


● H^+

RN 108357-04-6 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with 3-pyridinecarboxamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H^+

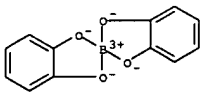
CM 2

CRN 108-47-4
CMF C7 H9 N

RN 108339-55-5 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-O,O']-, (T-4)-, hydrogen, compd. with piperazine (1:1) (9CI) (CA INDEX NAME)

CM 1

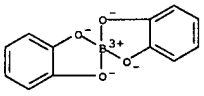
CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H^+

CM 2

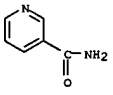
CRN 110-85-0
CMF C4 H10 N2



● H^+

CM 2

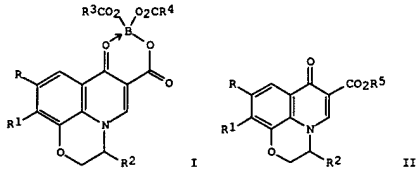
CRN 98-92-0
CMF C6 H6 N2 O



L12 ANSWER 93 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1986:553293 CAPLUS
DOCUMENT NUMBER: 105:153293
TITLE: Boron chelate compounds
PATENT ASSIGNEE(S): Daiichi Seiyaku Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 4 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 60075489	A2	19850427	JP 1983-184817	19831003
JP 04075239	B4	19921130		
PRIORITY APPLN. INFO.:			JP 1983-184817	19831003

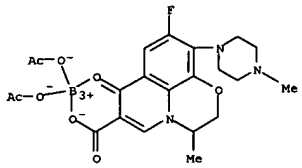
OTHER SOURCE(S): CASREACT 105:153293
GI



I II

AB Title chelates I (R, R1 = halo; R2 = H, alkyl; R3, R4 = aryl, alkyl, haloalkyl), intermediates for preparing antibacterial substances II [R1 = 4-(substituted)-1-piperazinyl; R5 = H], were prepared Thus, refluxing H3BO3, (EtCO)2O, and II (R = R1 = F; R2 = Me; R5 = Et) gave 95.2% I (R3 = R4 = Et) which was stirred with 4-methylpiperazine and neutralized to give 83.9% II (R1 = 4-methyl-1-piperazinyl; R5 = H).
IT **97746-90-2P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation and neutralization of)
RN 97746-90-2 CAPLUS
CN Boron, bis(acetato-O)[9-fluoro-2,3-dihydro-3-methyl-10-(4-methyl-1-piperazinyl)-7-oxo-7H-pyrido[1,2,3-de]-1,4-benzoxazine-6-carboxylato-O6,O7]-, (T-4)- (9CI) (CA INDEX NAME)

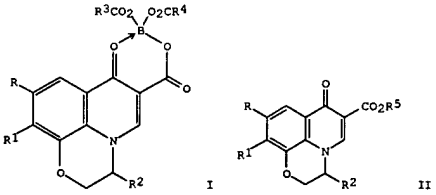
L12 ANSWER 93 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



L12 ANSWER 94 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1985:523491 CAPLUS
DOCUMENT NUMBER: 103:123491
TITLE: Oxazines
PATENT ASSIGNEE(S): Daiichi Seiyaku Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 6 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 60078986	A2	19850504	JP 1983-188138	19831007
JP 03072073	B4	19911115		
PRIORITY APPLN. INFO.:			JP 1983-188138	19831007

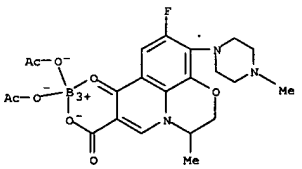
OTHER SOURCE(S): CASREACT 103:123491
GI



I II

AB Chelate dissociation of I [R = halo; R1 = (4-alkyl)-1-piperazinyl; R2 = H, alkyl; R3, R4 = aryl, alkyl, haloalkyl], prepared from I (R1 = halo) and (alkyl)piperazine, gave II having antibacterial activities. Thus, refluxing H3BO3, (EtCO)2O, and II (R = R1 = F; R2 = Me; R5 = Et) gave 95.2% I (R3 = R4 = Et), which was stirred with 4-methylpiperazine and neutralized to give 83.9% II (R1 = 4-methyl-1-piperazinyl; R5 = H).
IT **97746-90-2P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation and neutralization of)
RN 97746-90-2 CAPLUS
CN Boron, bis(acetato-O)[9-fluoro-2,3-dihydro-3-methyl-10-(4-methyl-1-piperazinyl)-7-oxo-7H-pyrido[1,2,3-de]-1,4-benzoxazine-6-carboxylato-O6,O7]-, (T-4)- (9CI) (CA INDEX NAME)

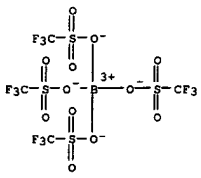
L12 ANSWER 94 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



L12 ANSWER 95 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1984:603144 CAPLUS
DOCUMENT NUMBER: 101:203144
TITLE: Contributions to the chemistry of boron, 148. Diamidoboron(1+) cations from bis(amino)borohalides by halide abstraction
AUTHOR(S): Noeth, Heinrich; Rasthofer, Bernhard; Weber, Siegfried
CORPORATE SOURCE: Inst. Anorg. Chem., Univ. Muenchen, Munich, D-8000/2, Fed. Rep. Ger.
SOURCE: Zeitschrift fuer Naturforschung, Teil B: Anorganische Chemie, Organische Chemie (1984), 39B(8), 1058-68
CODEN: ZNBAD2; ISSN: 0340-5087
DOCUMENT TYPE: Journal
LANGUAGE: German
AB Bis(amido)boron(1+) cations can be generated from bis(amino)boron halides and an appropriate halide acceptor. The ease of formation of the (Et2N)RB+ (R = 2,2,6,6-tetramethylpiperidinyl) cation from (Et2N)RBX (X = halide) was studied and found to increase with decreasing strength of the B-X bond. R2BF react with MCl3 (M = Al, Ga) or BX3 (X = Cl, Br, F) to give [R2B]MCl4 or [R2B]BX4, resp. [R2B]BBr4 reacts with AgO3SCF3 to yield [R2B]B(O3SCF3)4. R2BCl, obtained from [R2B]BCl4 in boiling pyridine, readily gives [R2B]M1Cl4 (M1 = metal). R2B(O3SCF3) is formed by reaction of R2BCl with AgO3SCF3.
IT **92816-40-5P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 92816-40-5 CAPLUS
CN Piperidinium, 2,2,6,6-tetramethyl-1-[(2,2,6,6-tetramethyl-1-piperidinyl)borylene]-, tetrakis(trifluoromethanesulfonato-O)borate(1-) (9CI) (CA INDEX NAME)

CM 1

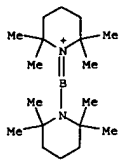
CRN 92816-39-2
CMF C4 B F12 O12 S4
CCI CCS



CM 2

CRN 92785-26-7
CMF C18 H36 B N2

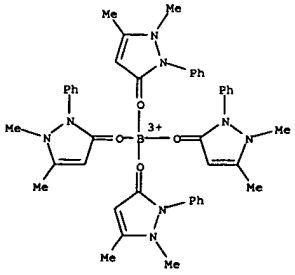
L12 ANSWER 95 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



L12 ANSWER 96 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1983:82717 CAPLUS
DOCUMENT NUMBER: 98:82717
TITLE: Acid-catalyzed electrophilic substitutions of 1,2-dihydro-1,5-dimethyl-2-phenyl-3H-pyrazol-3-one (phenazone)
AUTHOR(S): Akguen, Eyuep
CORPORATE SOURCE: Inst. Pharm. Lebensmittelchem., Univ. Wuerzburg, Wuerzburg, D-8700, Fed. Rep. Ger.
SOURCE: Chemiker-Zeitung (1982), 106(10), 371-3
CODEN: CMKZAT; ISSN: 0009-2894
DOCUMENT TYPE: Journal
LANGUAGE: German
AB [HRO]X (RO = 1,5-dimethyl-2-phenyl-3H-pyrazol-3-one; X = BF4, ClO4), [(RO)4B](SbCl6)3, and [(RO)4M](SbCl6)4 (M = Sn, Ti) were prepared and characterized by elemental anal. (except [ROH]X), 13C, 1H, 11B NMR and IR spectral measurements. These compds. are proposed as intermediates in the acid-catalyzed electrophilic substitution of RO.
IT **84663-13-8P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 84663-13-8 CAPLUS
CN Boron(3+), tetrakis(1,2-dihydro-1,5-dimethyl-2-phenyl-3H-pyrazol-3-one-O)-, (T-4)-, tris[(OC-6-11)-hexachloroantimonate(1-)] (9CI) (CA INDEX NAME)

CM 1

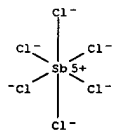
CRN 84663-12-7
CMF C44 H48 B N8 O4
CCI CCS



CM 2

CRN 17949-89-2
CMF Cl6 Sb
CCI CCS

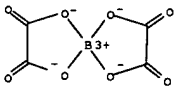
L12 ANSWER 96 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



L12 ANSWER 97 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1982:537689 CAPLUS
DOCUMENT NUMBER: 97:137689
TITLE: Boron complexes with dicarboxylic acids: bis(oxalato)borates and bis(malonato)borates
AUTHOR(S): Bessler, Eberhard; Weidlein, Johann
CORPORATE SOURCE: Dep. Quim., Univ. Brasilia, Brasilia, 70 910, Brazil
SOURCE: Zeitschrift fuer Naturforschung, Teil B: Anorganische Chemie, Organische Chemie (1982), 37B(8), 1020-5
CODEN: ZNBAD2; ISSN: 0340-5087
DOCUMENT TYPE: Journal
LANGUAGE: German
AB Several bis(oxalato)borates and bis(malonato)borates were prepared in heterogeneous reaction from boric acid, dicarboxylic acid and dicarboxylic acid salt by refluxing in benzene under continuous separation of H2O. The compds. were characterized by their IR and 11B NMR spectra. The vibrational spectra of the bis(oxalato)borate are discussed in detail.
IT **83145-84-0P 83145-92-0P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 83145-84-0 CAPLUS
CN Borate(1-), bis[ethanedioato(2-)-O,O']-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 83145-82-8
CMF C4 B O8 . H
CCI CCS



● H^+

CM 2

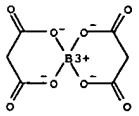
CRN 110-86-1
CMF C5 H5 N

RN 83145-92-0 CAPLUS

L12 ANSWER 97 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
CN Borate(1-), bis[propanedioato(2-)-O,O']-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 83145-90-8
CMF C6 H4 B O8 . H
CCI CCS



● H^+

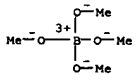
CM 2

CRN 110-86-1
CMF C5 H5 N

L12 ANSWER 98 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1982:113924 CAPLUS
DOCUMENT NUMBER: 96:113924
TITLE: Piperidinium tetramethoxyborate
AUTHOR(S): Alcock, Nathaniel W.; Hagger, Ruth M.; Harrison, W. David; Wallbridge, Malcolm G. H.
CORPORATE SOURCE: Dep. Chem. Mol. Sci., University of Warwick, Coventry, CV4 7AL, UK
SOURCE: Acta Crystallographica, Section B: Structural Crystallography and Crystal Chemistry (1982), B38(2), 676-7
CODEN: ACBCAR; ISSN: 0567-7408
DOCUMENT TYPE: Journal
LANGUAGE: English
AB The title salt is orthorhombic, space group Pnma, with a 13.325(3), b 9.603(2), and c 10.305(4) Å; Z = 4. The structure was solved by direct methods and refined to a final R = 0.051 for 494 reflections. Discrete cations and anions lying on mirror planes at b/2 are linked by H-bonds. Atomic coordinates are given.
IT **43160-43-6**
RL: PRP (Properties)
(structure of)
RN 43160-43-6 CAPLUS
CN Borate(1-), tetramethoxy-, hydrogen, compd. with piperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 49860-41-5
CMF C4 H12 B O4 . H
CCI CCS



● H^+

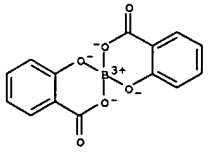
CM 2

CRN 110-89-4
CMF C5 H11 N

L12 ANSWER 99 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1979:72255 CAPLUS
DOCUMENT NUMBER: 90:72255
TITLE: The preparation and properties of some bis(salicylato)borate(III) salts with large cations
AUTHOR(S): Bassett, J.; Matthews, P. J.
CORPORATE SOURCE: Sch. Chem., Thames Polytech., London, UK
SOURCE: Journal of Inorganic and Nuclear Chemistry (1978), 40(6), 987-92
CODEN: JINCAO; ISSN: 0022-1902
DOCUMENT TYPE: Journal
LANGUAGE: English
AB Complexes of [BL2]- (HL = salicylic acid) with various cations (e.g., NH4, Bu4P, methylene blue) were prepared and characterized by anal. and IR, mass, and NMR spectral data. Analogous complexes with substituted salicylates were prepared and characterized similarly. The solvation of [NH4][BL2] (HL = 5-bromosalicylic acid) by MeCOR (R = Me, Et, iso-Bu) and Me2CHOH was studied.
IT **69030-71-3P 69030-85-9P 69030-91-7P 69030-97-3P 69030-98-4P 69030-99-5P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 69030-71-3 CAPLUS
CN Borate(1-), bis[2-hydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with piperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS



● H^+

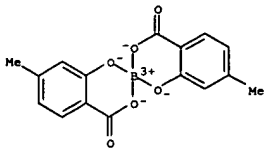
CM 2

CRN 110-89-4
CMF C5 H11 N

RN 69030-85-9 CAPLUS
CN Borate(1-), bis[2-hydroxy-4-methylbenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 69030-83-7
CMF C16 H12 B O6 . H
CCI CCS



● H+

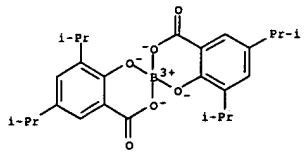
CM 2

CRN 110-86-1
CMF C5 H5 N

RN 69030-91-7 CAPLUS
CN Borate(1-), bis[2-hydroxy-3,5-bis(1-methylethyl)benzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 69030-89-3
CMF C26 H32 B O6 . H
CCI CCS



● H+

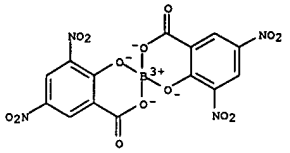
CM 2

CRN 110-86-1
CMF C5 H5 N

RN 69030-97-3 CAPLUS
CN Borate(1-), bis[2-hydroxy-3,5-dinitrobenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 4-ethylpyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 69030-96-2
CMF C14 H4 B N4 O14 . H
CCI CCS



● H+

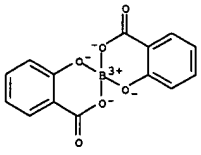
CM 2

CRN 536-75-4
CMF C7 H9 N

RN 69030-98-4 CAPLUS
CN Borate(1-), bis[2-(hydroxy-κO)benzoato(2-)-κO]-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS



● H+

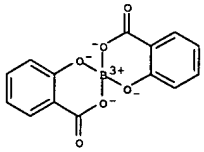
CM 2

CRN 110-86-1
CMF C5 H5 N

RN 69030-99-5 CAPLUS
CN Borate(1-), bis[2-hydroxybenzoato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with 2-methylpyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-97-1
CMF C14 H8 B O6 . H
CCI CCS

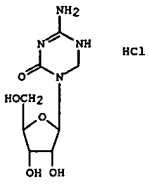


● H+

CM 2

CRN 109-06-8
CMF C6 H7 N

L12 ANSWER 100 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1977:433521 CAPLUS
DOCUMENT NUMBER: 87:33521
TITLE: Synthesis and antitumor activity of dihydro-5-azacytidine, a hydrolytically stable analog of 5-azacytidine
AUTHOR(S): Beisler, John A.; Abbasi, Mohamed M.; Kelley, James A.; Driscoll, John S.
CORPORATE SOURCE: Natl. Cancer Inst., NIH, Bethesda, MD, USA
SOURCE: Journal of Medicinal Chemistry (1977), 20(6), 806-12
CODEN: JMCMAR; ISSN: 0022-2623
DOCUMENT TYPE: Journal
LANGUAGE: English
GI



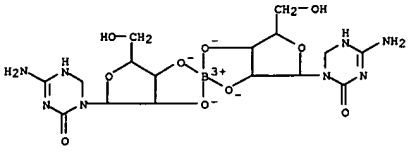
I

AB Borohydride reduction of 5-azacytidine [320-67-2] gave 5,6-dihydro-5-azacytidine-HCl (I) [62402-31-7] after acid hydrolysis of a boron-containing intermediate. Contrary to 5-azacytidine, which is easily hydrolyzed, I was completely stable at room temperature in aqueous solns. over a broad pH range for up to 3 weeks. I showed good activity in L1210 systems when administered i.p. or orally to mice. Although higher dose levels were necessary, I had approx. 80% of the antitumor efficacy shown by 5-azacytidine. Neither I nor 5-azacytidine showed a dependency on administration schedule. Cross-resistance between I and 5-azacytidine was demonstrated. I was superior to 5-azacytidine in therapeutic index and in its ability to cross the blood-brain barrier in sufficient quantity to be therapeutic against intracranially implanted L1210 cells. I may be a prodrug of 5-azacytidine.

IT **62769-14-6P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation and hydrolysis of)

RN 62769-14-6 CAPLUS
CN Borate(1-), bis[4-amino-1-β-D-ribofuranosyl-1,3,5-triazin-2(1H)-onato(2-)-O2',O3']-, sodium, (T-4)- (9CI) (CA INDEX NAME)

L12 ANSWER 100 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



● Na+

L12 ANSWER 101 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1974:5158 CAPLUS
DOCUMENT NUMBER: 80:5158
TITLE: Semipolar organoboron surfactants. I. Preparation of oil-soluble organoboron surfactants
AUTHOR(S): Hamanaka, Hiroyoshi
CORPORATE SOURCE: Surfactant Res. Lab., Toho Chem. Ind. Co., Tokyo, Japan
SOURCE: Yukagaku (1973), 22(8), 426-33
CODEN: YKGKAM; ISSN: 0513-398X
DOCUMENT TYPE: Journal
LANGUAGE: Japanese

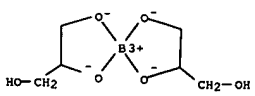
AB Glycerol [56-81-5] reacted with triethyl borate [150-46-9] at 130-50.deg. in 3 hr to give bisglycerol borate (I) [49625-59-4], which reacted further with fatty acids, e.g., caprylic acid [124-07-2], giving oil-soluble semipolar organoboron surfactants, such as bisglycerol borate dicaprylate [31973-04-3], with better surface-active properties than α-monoglycerides or their di-H borates.

IT **50258-91-8P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)

RN 50258-91-8 CAPLUS
CN Borate(1-), bis[1,2,3-propanetriolato(2-)-O1,O2]-, (T-4)-, hydrogen, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 49625-59-4
CMF C6 H12 B O6 . H
CCI CCS



● H+

CM 2

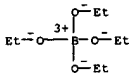
CRN 110-86-1
CMF C5 H5 N

L12 ANSWER 101 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

L12 ANSWER 102 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1973:532421 CAPLUS
DOCUMENT NUMBER: 79:132421
TITLE: Lewis acidity of trialkoxyboranes. Reinvestigation
AUTHOR(S): Wilson, John W.
CORPORATE SOURCE: Sch. Phys. Sci., New Univ. Ulster, Coleraine, UK
SOURCE: Journal of the Chemical Society, Dalton Transactions: Inorganic Chemistry (1972-1999) (1973), (16), 1628-30
CODEN: JCDTBI; ISSN: 0300-9246
DOCUMENT TYPE: Journal
LANGUAGE: English
AB B(OR)3 (R = Me, Et, Me2CH) formed onium salts containing [B(OR)4]- anions when reacted with piperidine, ethylenediamine, Et2NH, or cyclohexylamine in the presence of ROH, rather than Lewis acid-base adducts as previously reported (Ganbeau, J.; Link, R., 1951). The thermodn. stability of the salts is critically dependent on the crystal lattice energy. B(OPh)3 formed both L.B(OPh)3 and LH[B(OPh)4] (L = piperidine).
IT **43213-04-3P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 43213-04-3 CAPLUS
CN Borate(1-), tetraethoxy-, hydrogen, compd. with piperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 49861-34-9
CMF C8 H20 B O4 . H
CCI CCS



● H^+

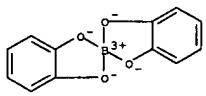
CM 2

CRN 110-89-4
CMF C5 H11 N

L12 ANSWER 103 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 1971:10679 CAPLUS
DOCUMENT NUMBER: 74:10679
TITLE: Combinations of boric acid with di- and triphenols
AUTHOR(S): Cismaru, D.; Cristescu, Virginia
CORPORATE SOURCE: Lab. Chim. Anorg., Fac. Farm., Bucharest, Rom.
SOURCE: Farmacia (Bucharest, Romania) (1970), 18(9), 531-7
CODEN: FRMBAZ; ISSN: 0014-8237
DOCUMENT TYPE: Journal
LANGUAGE: Romanian
GI For diagram(s), see printed CA Issue.
AB The complexes obtained by mixing boric acid with pyrocatechol or pyrogallol, and α-picoline or β-picoline, showed bactericidal activity against Salmonella, Shigella, Escherichia coli, and Proteus. As shown conductometrically, dipyrocatecholboric ions (I) and dipyrogallolboric ions were present in 0.01M solns.
IT **30776-62-6 30776-63-7 30776-64-8**
30776-65-9
RL: BAC (Biological activity or effector, except adverse); BSU (Biological
study, unclassified); BIOL (Biological study)
(bactericidal activity of)
RN 30776-62-6 CAPLUS
CN Borate(1-), bis[1,2-benzenediolato(2-)-κO,κO']-, (T-4)-, hydrogen, compd. with 2-methylpyridine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H^+

CM 2

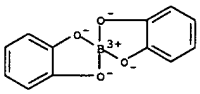
CRN 109-06-8
CMF C6 H7 N

RN 30776-63-7 CAPLUS
CN Borate(1-), bis[pyrocatecholato(2-)]-, hydrogen, compd. with 3-picoline

L12 ANSWER 103 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
(1:1) (8CI) (CA INDEX NAME)

CM 1

CRN 22450-98-2
CMF C12 H8 B O4 . H
CCI CCS



● H^+

CM 2

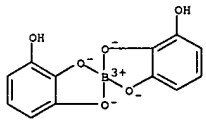
CRN 108-99-6
CMF C6 H7 N

RN 30776-64-8 CAPLUS
CN Borate(1-), bis[pyrogallolato(2-)]-, hydrogen, compd. with 2-picoline (8CI) (CA INDEX NAME)

CM 1

CRN 46944-61-0
CMF C12 H8 B O6 . H
CCI CCS

L12 ANSWER 103 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



● H^+

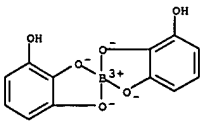
CM 2

CRN 109-06-8
CMF C6 H7 N

RN 30776-65-9 CAPLUS
CN Borate(1-), bis[pyrogallolato(2-)]-, hydrogen, compd. with 3-picoline (1:1) (8CI) (CA INDEX NAME)

CM 1

CRN 46944-61-0
CMF C12 H8 B O6 . H
CCI CCS



● H^+

CM 2

CRN 108-99-6
CMF C6 H7 N

L12 ANSWER 103 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

L12 ANSWER 104 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1969:14981 CAPLUS
DOCUMENT NUMBER: 70:14981
TITLE: Hydrolysis of trimethyl boric acid ester in the presence of organic bases
AUTHOR(S): Heller, G.
CORPORATE SOURCE: Freie Univ., Berlin, Fed. Rep. Ger.
SOURCE: Journal of Inorganic and Nuclear Chemistry (1968), 30(10), 2743-54
CODEN: JINCAO; ISSN: 0022-1902
DOCUMENT TYPE: Journal
LANGUAGE: German

AB Hydrolysis of trimethyl boric acid ester in the presence of organic bases in an organic solvent yields the following polyborates: the guanidinium salts [C(NH2)3]2[B3O3(OH)5], [C(NH2)3][B4O5(OH)5], [C(NH2)3]4[B5O6(OH)7]; the piperidinium salts [C5NH12][B4O4(OH)5], [C5NH12][B5O6(OH)4], the tetraalkylammonium salts [Me4N][B5O6(OH)4], [Et4N][B5O6(OH)4] (I), [(C3H7)4N][B7O6(OH)10] (II), [(C4H9)4N][B7O6(OH)10] (III), and the trialkylammonium borate [(C4H9)3NH][B7O6(OH)10] (IV). From aqueous solns. the evaporation of boric acid with an organic base results also in the formation of I, II, III, IV, and the guanidinium salt [C(NH2)3][B5O6(OH)4]. Only the compound [Me4N][B5O6(OH)4].4H2O containing crystal water forms an exception. All these polyborates undergo dissociation on dissoln., even in organic solvents like MeOH or Me2CO, to give lower mol. species. The structures of the new polyborates are discussed.

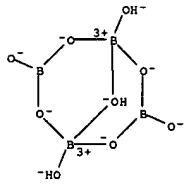
IT **12386-39-9**, Boric acid (H6B4O9), compound with piperidine (1:1)
12548-84-4, Boric acid (H5B5O10), compound with piperidine (1:1)
RL: PRP (Properties)
(spectrum of, ir)

RN 12386-39-9 CAPLUS
CN Borate(3-), μ-hydroxydihydroxybis[μ-[orthoborato(3-)-O:O']]di-, trihydrogen, compd. with piperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 12447-38-0
CMF B4 H3 O9 . 3 H
CCI CCS

L12 ANSWER 104 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



●3 H+

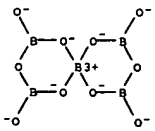
CM 2

CRN 110-89-4
CMF C5 H11 N

RN 12548-84-4 CAPLUS
CN Borate(5-), bis[μ-oxotetraoxodiborato(4-)]-, (T-4)-, pentahydrogen, compd. with piperidine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 92258-67-8
CMF B5 O10 . 5 H
CCI CCS



●5 H+

L12 ANSWER 104 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

CM 2

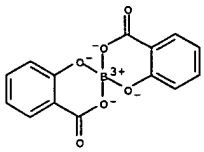
CRN 110-89-4
CMF C5 H11 N

L12 ANSWER 105 OF 105 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1968:56059 CAPLUS
DOCUMENT NUMBER: 68:56059
TITLE: Metal complexes of borosalicyclic acid and related substances
AUTHOR(S): Spacu, Petru; Gafiteanu, Mioara
CORPORATE SOURCE: Univ. Bucharest, Bucharest, Rom.
SOURCE: Analele Universitatii Bucuresti, Seria Stiintele Naturii (1966), 15(1), 33-49
CODEN: ABSNB3; ISSN: 0524-8302
DOCUMENT TYPE: Journal
LANGUAGE: Romanian
GI For diagram(s), see printed CA Issue.
AB Novel complexes were obtained between anions I, II, and III and the tetra-
and hexammine complexes of Co, Ni, and Cr. The uv and visible spectra were studied and conductometric measurements were made. The existence of the ions I, II, and III was confirmed via double exchange reactions. The complexes [Co(NH3)4(H2O)2]L3, [Co(NH3)6]L3, [Co(en)3]L3, [Co(pn)3]L3, [Co(dipy)3]L3, [Co(phen)3]L3, where L = [B(C7H4O3)2], dipy = dipyridyl, and phen = o-phenanthroline, were obtained from NH4[B(C7H4O3)2] and the resp. amine complexes in concentration aqueous solns., chilled with ice and under
constant agitation. The salts obtained with [Co(phen)3]Cl3 and [Co(dipy)3]Cl3 exhibited a greater stability and are more insol. than the other new complexes. The following were also prepared via double exchange
reactions: [Co(en)2Cl2]L, [Co(en)2(SCN)2]L, [Co(dipy)2Cl2]L, [Co(py)4Cl2]L. [Cr(NH3)6]L3, [Cr(en)3]L3, and [Cr(en)2(SCN)2]L were prepared as above. The preparation of [Co(en)3]L3 was difficult due to sensitivity to humidity and lack of stability at ambient conditions. [Ni(phen)3]L2 and [Ni(dipy)3]L2 were prepared via coupling with the Ni salts
and are stable in the absence of humidity. The salts of II were prepared from NH4[B(C6H4O2)2] and the resp. ammine complexes. [Co(dipy)3]X3 and [Co(phen)3]X3 were obtained, where X = B[(C6H4O2)2]. Other attempts to obtain complexes with various tetra-, penta, and hexammines were unsuccessful due to the highly dissociated state of the II ion in aqueous solution
The existence of the III ion in solution was evidenced for the 1st time by
the preparation of the complexes [Co(en)2Cl2]R and of [Co(en)2(SCN)2]R, which
are very stable, where R = [B(C7H5NO3)2].
IT **20149-68-2P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 20149-68-2 CAPLUS
CN Cobalt(1+), dichlorotetrakis(pyridine)-, bis[salicylato(2-)]borate(1-) (8CI) (CA INDEX NAME)

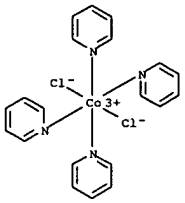
CM 1

CRN 38403-08-6
CMF C14 H8 B O6
CCI CCS



CM 2

CRN 18955-02-7
CMF C20 H20 Cl2 Co N4
CCI CCS



=> log y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	535.43	792.20

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-78.75	-78.75

STN INTERNATIONAL LOGOFF AT 09:17:38 ON 09 MAR 2006